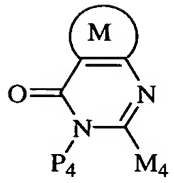
WHAT IS CLAIMED IS:

1. A compound of formula I:



I

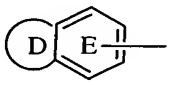
or a stereoisomer or pharmaceutically acceptable salt thereof, wherein;

ring M is a 5 or 6 membered aromatic or dihydro-aromatic ring consisting of: carbon atoms and 0-3 heteroatoms selected from O, $S(O)_p$, and N;

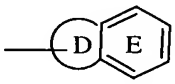
ring M is substituted with 0-3 R^{1a} and 0-1 carbonyl groups;

one of P_4 and M_4 is -Z-A-B and the other -G_1-G;

G is a group of formula IIa or IIb:



IIa



IIb

ring D, including the two atoms of Ring E to which it is attached, is a 5-6 membered ring consisting of: carbon atoms and 0-2 heteroatoms selected from the group consisting of N, O, and $S(O)_p$;

ring D is substituted with 0-2 R and has 0-3 ring double bonds;

E is selected from phenyl, pyridyl, pyrimidyl, pyrazinyl, and pyridazinyl, and is substituted with 1-2 R;

alternatively, ring D is absent and ring E is selected from phenyl, pyridyl, pyrimidyl, pyrazinyl, pyridazinyl, pyrrolyl, pyrazolyl, imidazolyl, isoxazolyl, oxazolyl, triazolyl, thienyl, and thiazolyl, and ring E is substituted with 1-2 R;

alternatively, ring D is absent and ring E is selected from phenyl, pyridyl, pyrimidyl, pyrazinyl, pyridazinyl, pyrrolyl, pyrazolyl, imidazolyl, isoxazolyl, oxazolyl, triazolyl, thienyl, and thiazolyl, and ring E is substituted with 1 R and with a 5-6 membered heterocycle consisting of: carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and $S(O)_p$, wherein the heterocycle is substituted with 0-1 carbonyls, 1-2 R, and 0-3 ring double bonds;

R is selected from H, $C_{1\text{-}4}$ alkyl, F, Cl, Br, I, OH, OCH_3, -OCH_2CH_3,

$-OCH(CH_3)_2$, $-OCH_2CH_2CH_3$, CN, $-C(=NR^8)NR^7R^9$, $-NHC(=NR^8)NR^7R^9$, $-NR^8CH(=NR^7)$, NH_2, $-NH(C_{1-3}$ alkyl), $-N(C_{1-3}$ alkyl)$_2$, $-C(=NH)NH_2$, $-CH_2NH_2$, $-CH_2NH(C_{1-3}$ alkyl), $-CH_2N(C_{1-3}$ alkyl)$_2$, $-CH_2CH_2NH_2$, $-CH_2CH_2NH(C_{1-3}$ alkyl), $-CH_2CH_2N(C_{1-3}$ alkyl)$_2$, $-(CR^8R^9)_tC(O)H$, $-(CR^8R^9)_tC(O)R^{2c}$, $-(CR^8R^9)_tNR^7R^8$, $-(CR^8R^9)_tC(O)NR^7R^8$, $-(CR^8R^9)_tNR^7C(O)R^7$, $-(CR^8R^9)_tOR^3$, $-(CR^8R^9)_tS(O)_pNR^7R^8$, $-(CR^8R^9)_tNR^7S(O)_pR^7$, $-(CR^8R^9)_tSR^3$, $-(CR^8R^9)_tS(O)R^3$, $-(CR^8R^9)_tS(O)_2R^3$, and $-OCF_3$, provided that $S(O)_pR^7$ forms other than $S(O)_2H$ or S(O)H;

alternatively, when 2 R groups are attached to adjacent atoms, they combine to form methylenedioxy or ethylenedioxy;

A is selected from: C_{3-10} carbocycle substituted with 0-2 R^4, and 5-12 membered heterocycle substituted with 0-2 R^4 and consisting of: carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and $S(O)_p$;

B is selected from: Y, X-Y, $-(CH_2)_{0-2}C(O)NR^2R^{2a}$, $-(CH_2)_{0-2}NR^2R^{2a}$, $-C(=NR^2)NR^2R^{2a}$, and $-NR^2C(=NR^2)NR^2R^{2a}$, provided that Z and B are attached to different atoms on A;

X is selected from $-(CR^2R^{2a})_{1-4}$-, $-CR^2(CR^2R^{2b})(CH_2)_t$-, -C(O)-, $-C(=NR^{1b})$-, $-CR^2(NR^{1b}R^2)$-, $-CR^2(OR^2)$-, $-CR^2(SR^2)$-, $-C(O)CR^2R^{2a}$-, $-CR^2R^{2a}C(O)$, -S-, -S(O)-, $-S(O)_2$-, $-SCR^2R^{2a}$-, $-S(O)CR^2R^{2a}$-, $-S(O)_2CR^2R^{2a}$-, $-CR^2R^{2a}S$-, $-CR^2R^{2a}S(O)$-, $-CR^2R^{2a}S(O)_2$-, $-S(O)_2NR^2$-, $-NR^2S(O)_2$-, $-NR^2S(O)_2CR^2R^{2a}$-, $-CR^2R^{2a}S(O)_2NR^2$-, $-NR^2S(O)_2NR^2$-, $-C(O)NR^2$-, $-NR^2C(O)$-, $-C(O)NR^2CR^2R^{2a}$-, $-NR^2C(O)CR^2R^{2a}$-, $-CR^2R^{2a}C(O)NR^2$-, $-CR^2R^{2a}NR^2C(O)$-, $-NR^2C(O)O$-, $-OC(O)NR^2$-, $-NR^2C(O)NR^2$-, $-NR^2$-, $-NR^2CR^2R^{2a}$-, $-CR^2R^{2a}NR^2$-, O, $-CR^2R^{2a}O$-, and $-OCR^2R^{2a}$-;

Y is selected from: C_{3-10} carbocycle substituted with 0-2 R^{4a}, and 5-10 membered heterocycle consisting of: carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and $S(O)_p$ substituted with 0-2 R^{4a};

G_1 is absent or is selected from $-(CR^3R^{3a})_{1-5}$-, $-(CR^3R^{3a})_{0-2}CR^3{=}CR^3(CR^3R^{3a})_{0-2}$-, $-(CR^3R^{3a})_{0-2}C{\equiv}C(CR^3R^{3a})_{0-2}$-, $-(CR^3R^{3a})_uC(O)(CR^3R^{3a})_w$-, $-(CR^3R^{3a})_uC(O)O(CR^3R^{3a})_w$-, $-(CR^3R^{3a})_uOC(O)(CR^3R^{3a})_w$-, $-(CR^3R^{3a})_uO(CR^3R^{3a})_w$-,

$-CR^3R^{3a})_uNR^{3b}(CR^3R^{3a})_w-$, $-(CR^3R^{3a})_uC(O)NR^{3b}(CR^3R^{3a})_w-$,
$-(CR^3R^{3a})_uNR^{3b}C(O)(CR^3R^{3a})_w-$, $-(CR^3R^{3a})_uOC(O)NR^{3b}(CR^3R^{3a})_w-$,
$-(CR^3R^{3a})_uNR^{3b}C(O)O(CR^3R^{3a})_w-$, $-(CR^3R^{3a})_uNR^{3b}C(O)NR^{3b}(CR^3R^{3a})_w-$,
$-(CR^3R^{3a})_uNR^{3^b}C(S)NR^{3b}(CR^3R^{3a})_w-$, $-(CR^3R^{3a})_uS(CR^3R^{3a})_w-$,
$-(CR^3R^{3a})_uS(O)(CR^3R^{3a})_w-$, $-(CR^3R^{3a})_uS(O)_2(CR^3R^{3a})_w-$,
$-(CR^3R^{3a})_uS(O)NR^{3b}(CR^3R^{3a})_w-$, $-(CR^3R^{3a})_uNR^{3b}S(O)_2(CR^3R^{3a})_w-$,
$-(CR^3R^{3a})_uS(O)_2NR^{3b}(CR^3R^{3a})_w-$, $-(CR^3R^{3a})_uNR^{3b}S(O)_2NR^{3b}(CR^3R^{3a})_w-$,
$-(CR^3R^{3a})_uNR^{3e}(CR^3R^{3a})_w-$, $-(CR^3R^{3a})_uC(O)(CR^3R^{3a})_uC(O)(CR^3R^{3a})_w-$,
$-(CR^3R^{3a})_uNR^{3b}(CR^3R^{3a})_uC(O)NR^{3b}(CR^3R^{3a})_w-$,
$-(CR^3R^{3a})_uNR^{3b}C(O)(CR^3R^{3a})_uC(O)(CR^3R^{3a})_w-$,
$-(CR^3R^{3a})_uC(O)(CR^3R^{3a})_uC(O)NR^{3b}(CR^3R^{3a})_w-$,
$-(CR^3R^{3a})_uNR^{3b}C(O)(CR^3R^{3a})_uC(O)NR^{3b}(CR^3R^{3a})_w-$,
$-(CR^3R^{3a})_uS(O)NR^{3b}C(O)(CR^3R^{3a})_w-$, $-(CR^3R^{3a})_uC(O)NR^{3b}S(O)_2(CR^3R^{3a})_w-$, and
$-(CR^3R^{3a})_uS(O)_2NR^{3b}C(O)NR^{3b}CR^3R^{3a})_w-$, wherein u + w total 0, 1, 2, 3, or 4,
provided that G_1 does not form a N-S, NCH_2N, NCH_2O, or NCH_2S bond with either group to which it is attached;

Z is selected from a bond, $-(CR^3R^{3a})_{1-4}-$, $-(CR^3R^{3a})_qO(CR^3R^{3a})_{q1}-$,
$-(CR^3R^{3a})_qNR^{3b}(CR^3R^{3a})_{q1}-$, $-(CR^3R^{3a})_qC(O)(CR^3R^{3a})_{q1}-$,
$-(CR^3R^{3a})_qC(O)O(CR^3R^{3a})_{q1}-$, $-(CR^3R^{3a})_qOC(O)(CR^3R^{3a})_{q1}-$,
$-(CR^3R^{3a})_qC(O)NR^{3b}(CR^3R^{3a})_{q1}-$, $-(CR^3R^{3a})_qNR^{3b}C(O)(CR^3R^{3a})_{q1}-$,
$-(CR^3R^{3a})_qOC(O)O(CR^3R^{3a})_{q1}-$, $-(CR^3R^{3a})_qOC(O)NR^{3b}(CR^3R^{3a})_{q1}-$,
$-(CR^3R^{3a})_qNR^{3b}C(O)O(CR^3R^{3a})_{q1}-$, $-(CR^3R^{3a})_qNR^{3b}C(O)NR^{3b}(CR^3R^{3a})_{q1}-$,
$-(CR^3R^{3a})_qC(O)(CR^3R^{3a})_qC(O)(CR^3R^{3a})_{q1}-$,
$-(CR^3R^{3a})_qNR^{3b}(CR^3R^{3a})_qC(O)NR^{3b}(CR^3R^{3a})_{q1}-$,
$-(CR^3R^{3a})_qNR^{3b}C(O)(CR^3R^{3a})_qC(O)(CR^3R^{3a})_{q1}-$,
$-(CR^3R^{3a})_qC(O)(CR^3R^{3a})_qC(O)NR^{3b}(CR^3R^{3a})_{q1}-$,
$-(CR^3R^{3a})_qNR^{3b}C(O)(CR^3R^{3a})_qC(O)NR^{3b}(CR^3R^{3a})_{q1}-$, $-(CR^3R^{3a})_qS(CR^3R^{3a})_{q1}-$,
$-(CR^3R^{3a})_qS(O)(CR^3R^{3a})_{q1}-$, $-(CR^3R^{3a})_qS(O)_2(CR^3R^{3a})_{q1}-$,
$-(CR^3R^{3a})_qSO_2NR^{3b}(CR^3R^{3a})_{q1}-$, $-(CR^3R^{3a})_qNR^{3b}SO_2(CR^3R^{3a})_{q1}-$,

$-(CR^3R^{3a})_qS(O)NR^{3b}C(O)(CR^3R^{3a})_{q1}-$, $-(CR^3R^{3a})_qC(O)NR^{3b}S(O)_2(CR^3R^{3a})_{q1}-$, and $-(CR^3R^{3a})_qNR^{3b}SO_2NR^{3b}(CR^3R^{3a})_{q1}-$, wherein q + q1 total 0, 1, 2, 3, or 4, provided that Z does not form a N-S, NCH_2N, NCH_2O, or NCH_2S bond with either group to which it is attached;

R^{1a}, at each occurrence, is selected from H, $-(CR^3R^{3a})_r-R^{1b}$, $-(CR^3R^{3a})_r-CR^3R^{1b}R^{1b}$, $-(CR^3R^{3a})_r-O-(CR^3R^{3a})_r-R^{1b}$, $-(CR^3R^{3a})_r-NR^2-(CR^3R^{3a})_r-R^{1b}$, $-(CR^3R^{3a})_r-S(O)_p-(CR^3R^{3a})_r-R^{1b}$, $-(CR^3R^{3a})_r-CO_2-(CR^3R^{3a})_r-R^{1b}$, $-(CR^3R^{3a})_r-C(O)NR^2-(CR^3R^{3a})_r-R^{1b}$, $-(CR^3R^{3a})_r-C(O)-(CR^3R^{3a})_r-R^{1b}$, $-C_{2-6}$ alkenylene-R^{1b}, $-C_{2-6}$ alkynylene-R^{1b}, and $-(CR^3R^{3a})_r-C(=NR^{1b})NR^3R^{1b}$, provided that R^{1a} forms other than an N-halo, N-S, O-O, or N-CN bond;

alternatively, when two R^{1a} groups are attached to adjacent atoms, together with the atoms to which they are attached, they form a 5-7 membered ring consisting of: carbon atoms and 0-2 heteroatoms selected from the group consisting of N, O, and $S(O)_p$, this ring being substituted with 0-2 R^{4b} and 0-3 ring double bonds;

R^{1b} is selected from H, C_{1-3} alkyl, F, Cl, Br, I, CN, NO_2, CHO, $-(CF_2)_rCF_3$, $-(CR^3R^{3a})_rOR^2$, $-NR^2R^{2a}$, $-C(O)R^{2b}$, $-CO_2R^{2b}$, $-OC(O)R^2$, $-CH(CH_2OR^2)_2$, $-(CF_2)_rCO_2R^{2a}$, $-S(O)_pR^{2b}$, $-NR^2(CH_2)_rOR^2$, $-C(=NR^{2c})NR^2R^{2a}$, $-NR^2C(O)R^{2b}$, $-NR^2C(O)NR^2R^{2a}$, $-NR^2C(O)_2R^{2a}$, $-OC(O)NR^2R^{2a}$, $-C(O)NR^2R^{2a}$, $-C(O)NR^2(CH_2)_rOR^2$, $-SO_2NR^2R^{2a}$, $-NR^2SO_2R^2$, $-C(O)NR^2SO_2R^2$, C_{3-6} carbocycle substituted with 0-2 R^{4b}, and 5-10 membered heterocycle substituted with 0-2 R^{4b} and consisting of carbon atoms and from 1-4 heteroatoms selected from the group consisting of N, O, and $S(O)_p$, provided that R^{1b} forms other than an O-O, N-halo, N-S, or N-CN bond and provided that $S(O)_pR^2$ forms other than $S(O)_2H$ or S(O)H;

R^2, at each occurrence, is selected from H, CF_3, C_{1-6} alkyl, benzyl, $-(CH_2)_r-C_{3-10}$ carbocycle substituted with 0-2 R^{4b}, and $-(CH_2)_r-$5-10 membered heterocycle substituted with 0-2 R^{4b} and consisting of: carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and $S(O)_p$;

R^{2a}, at each occurrence, is selected from H, CF_3, C_{1-6} alkyl, benzyl,

$-(CH_2)_r-C_{3-10}$ carbocycle substituted with 0-2 R^{4b}, and $-(CH_2)_r$-5-10 membered heterocycle substituted with 0-2 R^{4b} and consisting of: carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and $S(O)_p$;

alternatively, R^2 and R^{2a}, together with the nitrogen atom to which they are attached, combine to form a 5 or 6 membered saturated, partially saturated, or unsaturated ring substituted with 0-2 R^{4b} and consisting of: 0-1 additional heteroatoms selected from the group consisting of N, O, and $S(O)_p$;

R^{2b}, at each occurrence, is selected from CF_3, C_{1-4} alkoxy, C_{1-6} alkyl substituted with 0-2 R^{4b}, $-(CH_2)_r-C_{3-10}$ carbocycle substituted with 0-2 R^{4b}, and $-(CH_2)_r$-5-10 membered heterocycle substituted with 0-2 R^{4b} and consisting of: carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and $S(O)_p$;

R^{2c}, at each occurrence, is selected from CF_3, OH, C_{1-4} alkoxy, C_{1-6} alkyl, $-(CH_2)_r-C_{3-10}$ carbocycle substituted with 0-2 R^{4b}, and $-(CH_2)_r$-5-10 membered heterocycle substituted with 0-2 R^{4b} and consisting of carbon atoms and from 1-4 heteroatoms selected from the group consisting of N, O, and $S(O)_p$;

R^3, at each occurrence, is selected from H, CH_3, CH_2CH_3, $CH_2CH_2CH_3$, $CH(CH_3)_2$, $CH_2CH_2CH_2CH_3$, $CH_2CH(CH_3)_2$, $CH(CH_3)CH_2CH_3$, $C(CH_3)_3$, benzyl, and phenyl;

R^{3a}, at each occurrence, is selected from H, CH_3, CH_2CH_3, $CH_2CH_2CH_3$, $CH(CH_3)_2$, $CH_2CH_2CH_2CH_3$, $CH_2CH(CH_3)_2$, $CH(CH_3)CH_2CH_3$, $C(CH_3)_3$, benzyl, and phenyl;

R^{3b}, at each occurrence, is selected from H, C_{1-6} alkyl substituted with 0-2 R^{1a}, C_{2-6} alkenyl substituted with 0-2 R^{1a}, C_{2-6} alkynyl substituted with 0-2 R^{1a}, $-(C_{0-4}$ alkyl)-5-10 membered carbocycle substituted with 0-3 R^{1a}, and $-(C_{0-4}$ alkyl)- 5-10 membered heterocycle substituted with 0-3 R^{1a} and consisting of: carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and $S(O)_p$;

R^{3c}, at each occurrence, is selected from CH_3, CH_2CH_3, $CH_2CH_2CH_3$, $CH(CH_3)_2$, $CH_2CH_2CH_2CH_3$, $CH_2CH(CH_3)_2$, $CH(CH_3)CH_2CH_3$, $C(CH_3)_3$, benzyl, and phenyl;

R^{3d}, at each occurrence, is selected from H, CH_3, CH_2CH_3, $CH_2CH_2CH_3$, $CH(CH_3)_2$, $CH_2CH_2CH_2CH_3$, $CH_2CH(CH_3)_2$, $CH(CH_3)CH_2CH_3$, C_{1-4} alkyl-phenyl, and $C(=O)R^{3c}$;

R^{3e}, is selected from H, $-S(O)_2NHR^3$, $-C(O)R^3$, $-C(O)NHR^3$, $-C(O)OR^{3f}$, $-S(O)R^{3f}$, $-S(O)_2R^{3f}$, C_{1-6} alkyl substituted with 0-2 R^{1a}, C_{2-6} alkenyl substituted with 0-2 R^{1a}, C_{2-6} alkynyl substituted with 0-2 R^{1a}, $-(C_{0-4}$ alkyl)-5-10 membered carbocycle substituted with 0-3 R^{1a}, and $-(C_{0-4}$ alkyl)- 5-10 membered heterocycle substituted with 0-3 R^{1a} and consisting of: carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and $S(O)_p$;

R^{3f}, at each occurrence, is selected from: C_{1-6} alkyl substituted with 0-2 R^{1a}, C_{2-6} alkenyl substituted with 0-2 R^{1a}, C_{2-6} alkynyl substituted with 0-2 R^{1a}, $-(C_{0-4}$ alkyl)-5-10 membered carbocycle substituted with 0-3 R^{1a}, and $-(C_{0-4}$ alkyl)-5-10 membered heterocycle substituted with 0-3 R^{1a} and consisting of: carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and $S(O)_p$;

R^4, at each occurrence, is selected from H, =O, $(CR^3R^{3a})_rOR^2$, F, Cl, Br, I, C_{1-4} alkyl, $-(CR^3R^{3a})_rCN$, $-(CR^3R^{3a})_rNO_2$, $-(CR^3R^{3a})_rNR^2R^{2a}$, $-(CR^3R^{3a})_rC(O)R^{2c}$, $-(CR^3R^{3a})_rNR^2C(O)R^{2b}$, $-(CR^3R^{3a})_rC(O)NR^2R^{2a}$, $-(CR^3R^{3a})_rNR^2C(O)NR^2R^{2a}$, $-(CR^3R^{3a})_rC(=NR^2)NR^2R^{2a}$, $-(CR^3R^{3a})_rC(=NS(O)_2R^{5a})NR^2R^{2a}$, $-(CR^3R^{3a})_rNHC(=NR^2)NR^2R^{2a}$, $-(CR^3R^{3a})_rC(O)NHC(=NR^2)NR^2R^{2a}$, $-(CR^3R^{3a})_rSO_2NR^2R^{2a}$, $-(CR^3R^{3a})_rNR^2SO_2NR^2R^{2a}$, $-(CR^3R^{3a})_rNR^2SO_2$-C_{1-4} alkyl, $-(CR^3R^{3a})_rNR^2SO_2R^{5a}$, $-(CR^3R^{3a})_rS(O)_pR^{5a}$, $-(CR^3R^{3a})_r(CF_2)_rCF_3$, $-NHCH_2R^{1b}$, $-OCH_2R^{1b}$, $-SCH_2R^{1b}$, $-N(CH_2)_2(CH_2)_tR^{1b}$, $-O(CH_2)_2(CH_2)_tR^{1b}$, $-S(CH_2)_2(CH_2)_tR^{1b}$, $-(CR^3R^{3a})_r$-5-6 membered carbocycle substituted with 0-1 R^5, and a $-(CR^3R^{3a})_r$-5-6 membered heterocycle substituted with 0-1 R^5 and consisting of: carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and $S(O)_p$;

R^{4a}, at each occurrence, is selected from H, =O, -$(CR^3R^{3a})_rOR^2$, -$(CR^3R^{3a})_r$-F, -$(CR^3R^{3a})_r$-Br, -$(CR^3R^{3a})_r$-Cl, -$(CR^3R^{3a})_r$I, C_{1-4} alkyl, -$(CR^3R^{3a})_r$-CN, -$(CR^3R^{3a})_rNO_2$, -$(CR^3R^{3a})_rNR^2R^{2a}$, -$(CR^3R^{3a})_rC(O)R^{2c}$, -$(CR^3R^{3a})_rNR^2C(O)R^{2b}$, -$(CR^3R^{3a})_rC(O)NR^2R^{2a}$, -$(CR^3R^{3a})_rN{=}CHOR^3$, -$(CR^3R^{3a})_rC(O)NH(CH_2)_2NR^2R^{2a}$, -$(CR^3R^{3a})_rNR^2C(O)NR^2R^{2a}$, -$(CR^3R^{3a})_rNR^2C(O)OR^2$, -$(CR^3R^{3a})_rC(=NR^2)NR^2R^{2a}$, -$(CR^3R^{3a})_rNHC(=NR^2)NR^2R^{2a}$, -$(CR^3R^{3a})_rSO_2NR^2R^{2a}$, -$(CR^3R^{3a})_rNR^2SO_2NR^2R^{2a}$, -$(CR^3R^{3a})_rNR^2SO_2$-$C_{1-4}$ alkyl, -$(CR^3R^{3a})_rC(O)NHSO_2$-C_{1-4} alkyl, -$(CR^3R^{3a})_rNR^2SO_2R^5$, -$(CR^3R^{3a})_rS(O)_pR^5$, -$(CR^3R^{3a})_r(CF_2)_rCF_3$, -$(CR^3R^{3a})_r$-3-10 membered carbocycle substituted with 0-1 R^5, and a -$(CR^3R^{3a})_r$-3-10 membered heterocycle consisting of: carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and $S(O)_p$ and substituted with 0-1 R^5;

R^{4b}, at each occurrence, is selected from H, =O, -$(CR^3R^{3a})_rOR^3$, -$(CR^3R^{3a})_r$F,-$(CR^3R^{3a})_r$Cl, -$(CR^3R^{3a})_r$Br, -$(CR^3R^{3a})_r$I, C_{1-4} alkyl, -$(CR^3R^{3a})_r$CN, -$(CR^3R^{3a})_rNO_2$, -$(CR^3R^{3a})_rNR^3R^{3a}$, -$(CR^3R^{3a})_rC(O)R^3$, -$(CR^3R^{3a})_rC(O)OR^{3c}$, -$(CR^3R^{3a})_rNR^3C(O)R^{3a}$, -$(CR^3R^{3a})_rC(O)NR^3R^{3a}$, -$(CR^3R^{3a})_rNR^3C(O)NR^3R^{3a}$, -$(CR^3R^{3a})_rC(=NR^3)NR^3R^{3a}$, -$(CR^3R^{3a})_rNR^3C(=NR^3)NR^3R^{3a}$, -$(CR^3R^{3a})_rSO_2NR^3R^{3a}$, -$(CR^3R^{3a})_rNR^3SO_2NR^3R^{3a}$, -$(CR^3R^{3a})_rNR^3SO_2$-$C_{1-4}$ alkyl, -$(CR^3R^{3a})_rNR^3SO_2CF_3$, -$(CR^3R^{3a})_rNR^3SO_2$-phenyl, -$(CR^3R^{3a})_rS(O)_pCF_3$, -$(CR^3R^{3a})_rS(O)_p$-$C_{1-4}$ alkyl, -$(CR^3R^{3a})_rS(O)_p$-phenyl, and -$(CR^3R^{3a})_r(CF_2)_rCF_3$;

R^5, at each occurrence, is selected from H, C_{1-6} alkyl, =O, -$(CH_2)_rOR^3$, F, Cl, Br, I, -CN, NO_2, -$(CH_2)_rNR^3R^{3a}$, -$(CH_2)_rC(O)R^3$, -$(CH_2)_rC(O)OR^{3c}$, -$NR^3C(O)R^{3a}$, -$C(O)NR^3R^{3a}$, -$NR^3C(O)NR^3R^{3a}$, -$CH(=NOR^{3d})$, -$C(=NR^3)NR^3R^{3a}$, -$NR^3C(=NR^3)NR^3R^{3a}$, -$SO_2NR^3R^{3a}$, -$NR^3SO_2NR^3R^{3a}$, -NR^3SO_2-C_{1-4} alkyl, -$NR^3SO_2CF_3$, -NR^3SO_2-phenyl, -$S(O)_pCF_3$, -$S(O)_p$-C_{1-4} alkyl, -$S(O)_p$-phenyl, -$(CF_2)_rCF_3$, phenyl substituted with 0-2 R^6, naphthyl substituted with 0-2 R^6, and benzyl substituted with 0-2 R^6;

R^{5a}, at each occurrence, is selected from C_{1-6} alkyl, -$(CH_2)_rOR^3$,

$-(CH_2)_rNR^3R^{3a}$, $-(CH_2)_rC(O)R^3$, $-(CH_2)_rC(O)OR^{3c}$, $-(CH_2)_rNR^3C(O)R^{3a}$, $-(CH_2)_rC(O)NR^3R^{3a}$, $-(CF_2)_rCF_3$, phenyl substituted with 0-2 R^6, naphthyl substituted with 0-2 R^6, and benzyl substituted with 0-2 R^6, provided that R^{5a} does not form a S-N or $S(O)_p$-C(O) bond;

R^6, at each occurrence, is selected from H, OH, $-(CH_2)_rOR^2$, Fl, Cl, Br, I, C_{1-4} alkyl, -CN, NO_2, $-(CH_2)_rNR^2R^{2a}$, $-(CH_2)_rC(O)R^{2b}$, $-NR^2C(O)R^{2b}$, $-NR^2C(O)NR^2R^{2a}$, $-C(=NH)NH_2$, $-NHC(=NH)NH_2$, $-SO_2NR^2R^{2a}$, $-NR^2SO_2NR^2R^{2a}$, and $-NR^2SO_2C_{1-4}$ alkyl;

R^7, at each occurrence, is selected from H, OH, C_{1-6} alkyl, C_{1-6} alkyl-C(O)-, C_{1-6} alkyl-O-, $(CH_2)_n$-phenyl, C_{1-4} alkyl-OC(O)-, C_{6-10} aryl-O-, C_{6-10} aryl-OC(O)-, C_{6-10} aryl-CH_2-C(O)-, C_{1-4} alkyl-C(O)O-C_{1-4} alkyl-OC(O)-, C_{6-10} aryl-C(O)O-C_{1-4} alkyl-OC(O)-, C_{1-6} alkyl-NH_2-C(O)-, phenyl-NH_2-C(O)-, and phenyl C_{1-4} alkyl-C(O)-;

R^8, at each occurrence, is selected from H, C_{1-6} alkyl, and - $(CH_2)_n$-phenyl;

alternatively, R^7 and R^8, when attached to the same nitrogen, combine to form a 5-10 membered heterocyclic ring consisting of carbon atoms and 0-2 additional heteroatoms selected from the group consisting of N, O, and $S(O)_p$;

R^9, at each occurrence, is selected from H, C_{1-6} alkyl, and $-(CH_2)_n$-phenyl;

n, at each occurrence, is selected from 0, 1, 2, and 3;

p, at each occurrence, is selected from 0, 1, and 2;

r, at each occurrence, is selected from 0, 1, 2, 3, 4, 5, and 6; and

t, at each occurrence, is selected from 0, 1, 2, and 3;

provided that:

(a) when Z and G_1 are absent, A is phenyl or pyridyl, G is phenyl, pyridyl, or thienyl, at least one R is other than a substituted or unsubstituted group selected from amidino, guanidino, guanidine-methyl, iminoamino, iminoamino-methyl, amino, amino-methyl, and pyridyl, then B is other than cycloalkyl, $(CH_2)_{0-2}C(O)NR^2R^{2a}$, or $(CH_2)_{0-2}NR^2R^{2a}$, wherein substituted includes being cyclized with an additional heteroatom being optionally present;

(b) when Z and G_1 are absent, G is phenyl or pyridyl, and A is phenyl, pyridyl, furanyl, or thienyl, then B is other than a substituted or unsubstituted group selected from amidino, guanidino, guanidine-methyl, iminoamino, iminoamino-methyl, amino, amino-methyl, aminosulfonyl-phenyl, and pyridyl, wherein substituted includes being cyclized with an additional heteroatom being optionally present; and

(c) when $G\text{-}G_1$ is hydroxy-phenyl or alkoxy-phenyl, then B is other than acyclic or cyclic-amino-alkoxy.

2. A compound according to Claim 1, wherein:

ring M is selected from phenyl, pyrrole, furan, thiophene, pyrazole, imidazole, isoxazole, oxazole, isothiazole, thiazole, 1,2,3-triazole, 1,2,3-oxadiazole, 1,2,3-thiadiazole, pyridine, pyrimidine, pyridazine, pyrazine, 1,2,3-triazine, 1,2,4-triazine, and 1,2,3,4-tetrazine;

ring M is substituted with 0-3 R^{1a};

one of P_4 and M_4 is –Z-A-B and the other $-G_1$-G;

G is a group of formula IIa or IIb:

IIa IIb

ring D, including the two atoms of Ring E to which it is attached, is a 5-6 membered ring consisting of: carbon atoms and 0-2 heteroatoms selected from the group consisting of N, O, and $S(O)_p$;

ring D is substituted with 0-2 R and has 0-3 ring double bonds;

E is selected from phenyl, pyridyl, pyrimidyl, pyrazinyl, and pyridazinyl, and is substituted with 1-2 R;

alternatively, ring D is absent, and ring E is selected from phenyl, pyridyl, pyrimidyl, and thienyl, and ring E is substituted with 1-2 R;

alternatively, ring D is absent, ring E is selected from phenyl, pyridyl, and thienyl, and ring E is substituted with a 5 membered heterocycle consisting of: carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and $S(O)_p$,

wherein the heterocycle is substituted with 0-1 carbonyls, 1-2 R, and 0-3 ring double bonds;

R is selected from H, C_{1-4} alkyl, F, Cl, OH, OCH_3, $-OCH_2CH_3$, $-OCH(CH_3)_2$, CN, $-C(=NH)NH_2$, NH_2, $-NH(C_{1-3}$ alkyl), $-N(C_{1-3}$ alkyl$)_2$, $-C(=NH)NH_2$, $-CH_2NH_2$, $-CH_2NH(C_{1-3}$ alkyl), $-CH_2N(C_{1-3}$ alkyl$)_2$, $-(CR^8R^9)_tNR^7R^8$, $-C(O)NR^7R^8$, $-CH_2C(O)NR^7R^8$, $-S(O)_pNR^7R^8$, $-CH_2S(O)_pNR^7R^8$, and $-OCF_3$;

alternatively, when 2 R groups are attached to adjacent atoms, they combine to form methylenedioxy or ethylenedioxy;

A is selected from one of the following rings and is substituted with 0-2 R^4;

phenyl, piperidinyl, piperazinyl, pyridyl, pyrimidyl, furanyl, morpholinyl, thiophenyl, pyrrolyl, pyrrolidinyl, oxazolyl, isoxazolyl, thiazolyl, isothiazolyl, pyrazolyl, imidazolyl, oxadiazolyl, thiadiazolyl, triazolyl, 1,2,3-oxadiazolyl, 1,2,4-oxadiazolyl, 1,2,5-oxadiazolyl, 1,3,4-oxadiazolyl, 1,2,3-thiadiazolyl, 1,2,4-thiadiazolyl, 1,2,5-thiadiazolyl, 1,3,4-thiadiazolyl, 1,2,3-triazolyl, 1,2,4-triazolyl, 1,2,5-triazolyl, 1,3,4-triazolyl, benzofuranyl, benzothiofuranyl, indolyl, benzimidazolyl, benzoxazolyl, benzthiazolyl, indazolyl, benzisoxazolyl, benzisothiazolyl, and isoindazolyl;

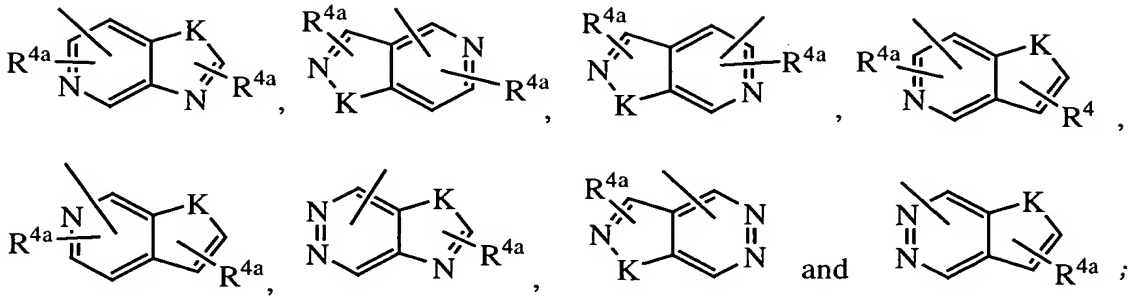
B is selected from Y, X-Y, $-CH_2NR^2R^{2a}$, and $-CH_2CH_2NR^2R^{2a}$;

X is selected from $-(CR^2R^{2a})_{1-4}$-, -C(O)-, $-C(=NR^{1b})$-, $-CR^2(NR^{1b}R^2)$-, $-C(O)CR^2R^{2a}$-, $-CR^2R^{2a}C(O)$, $-C(O)NR^2$-, $-NR^2C(O)$-, $-C(O)NR^2CR^2R^{2a}$-, $-NR^2C(O)CR^2R^{2a}$-, $-CR^2R^{2a}C(O)NR^2$-, $-CR^2R^{2a}NR^2C(O)$-, $-NR^2C(O)NR^2$-, $-NR^2$-, $-NR^2CR^2R^{2a}$-, $-CR^2R^{2a}NR^2$-, O, $-CR^2R^{2a}O$-, and $-OCR^2R^{2a}$-;

Y is selected from one of the following rings and is substituted with 0-2 R^{4a};

cyclopropyl, cyclopentyl, cyclohexyl, phenyl, piperidinyl, piperazinyl, pyridyl, pyrimidyl, furanyl, morpholinyl, thiophenyl, pyrrolyl, pyrrolidinyl, oxazolyl, isoxazolyl, isoxazolinyl, thiazolyl, isothiazolyl, pyrazolyl, imidazolyl, 1,2,3-oxadiazolyl, 1,2,4-oxadiazolyl, 1,2,5-oxadiazolyl, 1,3,4-oxadiazolyl, 1,2,3-thiadiazolyl, 1,2,4-thiadiazolyl, 1,2,5-thiadiazolyl, 1,3,4-thiadiazolyl, 1,2,3-triazolyl, 1,2,4-triazolyl, 1,2,5-triazolyl, 1,3,4-triazolyl, benzofuranyl, benzothiofuranyl, indolyl, benzimidazolyl, benzoxazolyl, benzthiazolyl, indazolyl, benzisoxazolyl, benzisothiazolyl, and isoindazolyl;

alternatively, Y is selected from the following bicyclic heteroaryl ring systems:



K is selected from O, S, NH, and N;

G_1 is absent or is selected from $-(CR^3R^{3a})_{1-3}-$, $-(CR^3R^{3a})_uC(O)(CR^3R^{3a})_w-$, $-(CR^3R^{3a})_uO(CR^3R^{3a})_w-$, $-(CR^3R^{3a})_uNR^{3b}(CR^3R^{3a})_w-$, $-(CR^3R^{3a})_uC(O)NR^{3b}(CR^3R^{3a})_w-$, $-(CR^3R^{3a})_uNR^{3b}C(O)(CR^3R^{3a})_w-$, $-(CR^3R^{3a})_uNR^{3b}C(O)(CR^3R^{3a})_uC(O)NR^{3b}(CR^3R^{3a})_w-$, $-(CR^3R^{3a})_uS(CR^3R^{3a})_w-$, $-(CR^3R^{3a})_uS(O)(CR^3R^{3a})_w-$, $-(CR^3R^{3a})_uS(O)_2(CR^3R^{3a})_w-$, $-(CR^3R^{3a})_uS(O)NR^{3b}(CR^3R^{3a})_w-$, $-(CR^3R^{3a})_uNR^{3b}S(O)_2(CR^3R^{3a})_w-$, and $-(CR^3R^{3a})_uS(O)_2NR^{3b}(CR^3R^{3a})_w-$, wherein u + w total 0, 1, or 2, provided that G_1 does not form a N-S, NCH_2N, NCH_2O, or NCH_2S bond with either group to which it is attached;

Z is selected from a bond, CH_2, CH_2CH_2, CH_2O, OCH_2, C(O), NH, CH_2NH, $NHCH_2$, $CH_2C(O)$, $C(O)CH_2$, C(O)NH, NHC(O), $NHC(O)CH_2C(O)NH$, $S(O)_2$, $CH_2S(O)_2$, $S(O)_2(CH_2)$, SO_2NH, and $NHSO_2$, provided that Z does not form a N-S, NCH_2N, NCH_2O, or NCH_2S bond with either group to which it is attached;

R^{1a}, at each occurrence, is selected from H, $-(CH_2)_r-R^{1b}$, $-(CH(CH_3))_r-R^{1b}$, $-(C(CH_3)_2)_r-R^{1b}$, $-O-(CR^3R^{3a})_r-R^{1b}$, $-NR^2-(CR^3R^{3a})_r-R^{1b}$, and $-S-(CR^3R^{3a})_r-R^{1b}$, provided that R^{1a} forms other than an N-halo, N-S, O-O, or N-CN bond;

alternatively, when two R^{1a} groups are attached to adjacent atoms, together with the atoms to which they are attached they form a 5-7 membered ring consisting of: carbon atoms and 0-2 heteroatoms selected from the group consisting of N, O, and $S(O)_p$, this ring being substituted with 0-2 R^{4b} and 0-3 ring double bonds;

R^{1b} is selected from H, CH_3, CH_2CH_3, $CH_2CH_2CH_3$, $CH(CH_3)_2$, F, Cl, Br, I, CN, CHO, CF_3, OR^2, $-NR^2R^{2a}$, $-C(O)R^{2b}$, $-CO_2R^{2b}$, $-OC(O)R^2$, $-CO_2R^{2a}$,

$-S(O)_pR^{2b}$, $-NR^2(CH_2)_rOR^2$, $-NR^2C(O)R^{2b}$, $-NR^2C(O)NHR^2$, $-NR^2C(O)_2R^{2a}$, $-OC(O)NR^2R^{2a}$, $-C(O)NR^2R^{2a}$, $-C(O)NR^2(CH_2)_rOR^2$, $-SO_2NR^2R^{2a}$, $-NR^2SO_2R^2$, C_{5-6} carbocycle substituted with 0-2 R^{4b}, and 5-6 membered heterocycle substituted with 0-2 R^{4b} and consisting of carbon atoms and from 1-4 heteroatoms selected from the group consisting of N, O, and $S(O)_p$, provided that R^{1b} forms other than an O-O, N-halo, N-S, or N-CN bond and provided that $S(O)_pR^2$ forms other than $S(O)_2H$ or S(O)H;

R^2, at each occurrence, is selected from H, CF_3, CH_3, CH_2CH_3, $CH_2CH_2CH_3$, $CH(CH_3)_2$, $CH_2CH_2CH_2CH_3$, $CH_2CH(CH_3)_2$, $CH(CH_3)CH_2CH_3$, $C(CH_3)_3$, benzyl, C_{3-6} carbocycle substituted with 0-2 R^{4b}, C_{3-6} carbocycle-CH_2- substituted with 0-2 R^{4b}, and 5-6 membered heterocycle substituted with 0-2 R^{4b} and consisting of: carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and $S(O)_p$;

R^{2a}, at each occurrence, is selected from H, CF_3, CH_3, CH_2CH_3, $CH_2CH_2CH_3$, $CH(CH_3)_2$, $CH_2CH_2CH_2CH_3$, $CH_2CH(CH_3)_2$, $CH(CH_3)CH_2CH_3$, $C(CH_3)_3$, benzyl, C_{5-6} carbocycle substituted with 0-2 R^{4b}, and 5-6 membered heterocycle substituted with 0-2 R^{4b} and consisting of: carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and $S(O)_p$;

alternatively, R^2 and R^{2a}, together with the nitrogen atom to which they are attached, combine to form a 5 or 6 membered saturated, partially saturated, or unsaturated ring substituted with 0-2 R^{4b} and consisting of: 0-1 additional heteroatoms selected from the group consisting of N, O, and $S(O)_p$;

R^{2b}, at each occurrence, is selected from CF_3, C_{1-4} alkoxy, CH_3, CH_2CH_3, $CH_2CH_2CH_3$, $CH(CH_3)_2$, $CH_2CH_2CH_2CH_3$, $CH_2CH(CH_3)_2$, $CH(CH_3)CH_2CH_3$, $C(CH_3)_3$, benzyl, C_{5-6} carbocycle substituted with 0-2 R^{4b}, and 5-6 membered heterocycle substituted with 0-2 R^{4b} and consisting of: carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and $S(O)_p$;

R^{2c}, at each occurrence, is selected from CF_3, OH, C_{1-4} alkoxy, CH_3, CH_2CH_3, $CH_2CH_2CH_3$, $CH(CH_3)_2$, $CH_2CH_2CH_2CH_3$, $CH_2CH(CH_3)_2$, $CH(CH_3)CH_2CH_3$, $C(CH_3)_3$, benzyl, C_{5-6} carbocycle substituted with 0-2 R^{4b}, and

5-6 membered heterocycle substituted with 0-2 R^{4b} and consisting of carbon atoms and from 1-4 heteroatoms selected from the group consisting of N, O, and $S(O)_p$;

R^3, at each occurrence, is selected from H, CH_3, CH_2CH_3, $CH_2CH_2CH_3$, $CH(CH_3)_2$, benzyl, and phenyl;

R^{3a}, at each occurrence, is selected from H, CH_3, CH_2CH_3, $CH_2CH_2CH_3$, $CH(CH_3)_2$, benzyl, and phenyl;

R^{3c}, at each occurrence, is selected from CH_3, CH_2CH_3, $CH_2CH_2CH_3$, $CH(CH_3)_2$, benzyl, and phenyl;

R^{3d}, at each occurrence, is selected from H, CH_3, CH_2CH_3, $CH_2CH_2CH_3$, $CH(CH_3)_2$, CH_2-phenyl, CH_2CH_2-phenyl, and $C(=O)R^{3c}$;

R^4, at each occurrence, is selected from H, =O, OR^2, $-CH_2OR^2$, $-(CH_2)_2OR^2$, F, Cl, Br, I, C_{1-4} alkyl, CN, NO_2, $-NR^2R^{2a}$, $-CH_2NR^2R^{2a}$, $-(CH_2)_2NR^2R^{2a}$, $-C(O)R^{2c}$, $-NR^2C(O)R^{2b}$, $-C(O)NR^2R^{2a}$, $-SO_2NR^2R^{2a}$, $-S(O)_pR^{5a}$, CF_3, CF_2CF_3, 5-6 membered carbocycle substituted with 0-1 R^5, and a 5-6 membered heterocycle substituted with 0-1 R^5 and consisting of: carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and $S(O)_p$;

R^{4a}, at each occurrence, is selected from H, =O, $-(CR^3R^{3a})_rOR^2$, $-(CR^3R^{3a})_r$-F, $-(CR^3R^{3a})_r$-Br, $-(CR^3R^{3a})_r$-Cl, C_{1-4} alkyl, $-(CR^3R^{3a})_r$-CN, $-(CR^3R^{3a})_rNO_2$, $-(CR^3R^{3a})_rNR^2R^{2a}$, $-(CR^3R^{3a})_rC(O)R^{2c}$, $-(CR^3R^{3a})_rNR^2C(O)R^{2b}$, $-(CR^3R^{3a})_rC(O)NR^2R^{2a}$, $-(CR^3R^{3a})_rSO_2NR^2R^{2a}$, $-(CR^3R^{3a})_rNR^2SO_2NR^2R^{2a}$, $-(CR^3R^{3a})_rNR^2SO_2$-C_{1-4} alkyl, $-(CR^3R^{3a})_rC(O)NHSO_2$-$C_{1-4}$ alkyl, $-(CR^3R^{3a})_rNR^2SO_2R^5$, $-(CR^3R^{3a})_rS(O)_pR^5$, $-(CR^3R^{3a})_r(CF_2)_rCF_3$, phenyl substituted with 0-1 R^5, and a 5 membered aromatic heterocycle consisting of: carbon atoms and 1-3 heteroatoms selected from the group consisting of N, O, and $S(O)_p$ substituted with 0-1 R^5;

R^{4b}, at each occurrence, is selected from H, =O, OR^3, CH_2OR^3, F, Cl, CH_3, CH_2CH_3, $CH_2CH_2CH_3$, $CH(CH_3)_2$, $CH_2CH_2CH_2CH_3$, $CH_2CH(CH_3)_2$, $CH(CH_3)CH_2CH_3$, $C(CH_3)_3$, CN, NO_2, $-NR^3R^{3a}$, $-CH_2NR^3R^{3a}$, $-C(O)R^3$, $-CH_2$-$C(O)R^3$, $-C(O)OR^{3c}$, $-CH_2C(O)OR^{3c}$, $-NR^3C(O)R^{3a}$, $-CH_2NR^3C(O)R^{3a}$,

$-C(O)NR^3R^{3a}$, $-CH_2C(O)NR^3R^{3a}$, $-NR^3C(O)NR^3R^{3a}$, $-CH_2NR^3C(O)NR^3R^{3a}$, $-C(=NR^3)NR^3R^{3a}$, $-CH_2C(=NR^3)NR^3R^{3a}$, $-NR^3C(=NR^3)NR^3R^{3a}$, $-CH_2NR^3C(=NR^3)NR^3R^{3a}$, $-SO_2NR^3R^{3a}$, $-CH_2SO_2NR^3R^{3a}$, $-NR^3SO_2NR^3R^{3a}$, $-CH_2NR^3SO_2NR^3R^{3a}$, $-NR^3SO_2$-C_{1-4} alkyl, $-CH_2NR^3SO_2$-C_{1-4} alkyl, $-NR^3SO_2CF_3$, $-CH_2NR^3SO_2CF_3$, $-NR^3SO_2$-phenyl, $-CH_2NR^3SO_2$-phenyl, $-S(O)_pCF_3$, $-CH_2S(O)_pCF_3$, $-S(O)_p$-C_{1-4} alkyl, $-CH_2S(O)_p$-C_{1-4} alkyl, $-S(O)_p$-phenyl, $-CH_2S(O)_p$-phenyl, CF_3, and CH_2-CF_3;

R^5, at each occurrence, is selected from H, =O, CH_3, CH_2CH_3, $CH_2CH_2CH_3$, $CH(CH_3)_2$, $CH_2CH_2CH_2CH_3$, $CH_2CH(CH_3)_2$, $CH(CH_3)CH_2CH_3$, $C(CH_3)_3$, OR^3, $-CH_2OR^3$, F, Cl, CN, NO_2, $-NR^3R^{3a}$, $-CH_2NR^3R^{3a}$, $-C(O)R^3$, $-CH_2C(O)R^3$, $-C(O)OR^{3c}$, $-CH_2C(O)OR^{3c}$, $-NR^3C(O)R^{3a}$, $-C(O)NR^3R^{3a}$, $-NR^3C(O)NR^3R^{3a}$, $-CH(=NOR^{3d})$, $-C(=NR^3)NR^3R^{3a}$, $-NR^3C(=NR^3)NR^3R^{3a}$, $-SO_2NR^3R^{3a}$, $-NR^3SO_2NR^3R^{3a}$, $-NR^3SO_2$-C_{1-4} alkyl, $-NR^3SO_2CF_3$, $-NR^3SO_2$-phenyl, $-S(O)_pCF_3$, $-S(O)_p$-C_{1-4} alkyl, $-S(O)_p$-phenyl, CF_3, phenyl substituted with 0-2 R^6, naphthyl substituted with 0-2 R^6, and benzyl substituted with 0-2 R^6;

R^6, at each occurrence, is selected from H, OH, OR^2, F, Cl, CH_3, CH_2CH_3, $CH_2CH_2CH_3$, $CH(CH_3)_2$, $CH_2CH_2CH_2CH_3$, $CH_2CH(CH_3)_2$, $CH(CH_3)CH_2CH_3$, $C(CH_3)_3$, CN, NO_2, $-NR^2R^{2a}$, $-CH_2NR^2R^{2a}$, $-C(O)R^{2b}$, $-CH_2C(O)R^{2b}$, $-NR^2C(O)R^{2b}$, $-NR^2C(O)NR^2R^{2a}$, $-C(=NH)NH_2$, $-NHC(=NH)NH_2$, $-SO_2NR^2R^{2a}$, $-NR^2SO_2NR^2R^{2a}$, and $-NR^2SO_2C_{1-4}$ alkyl; and

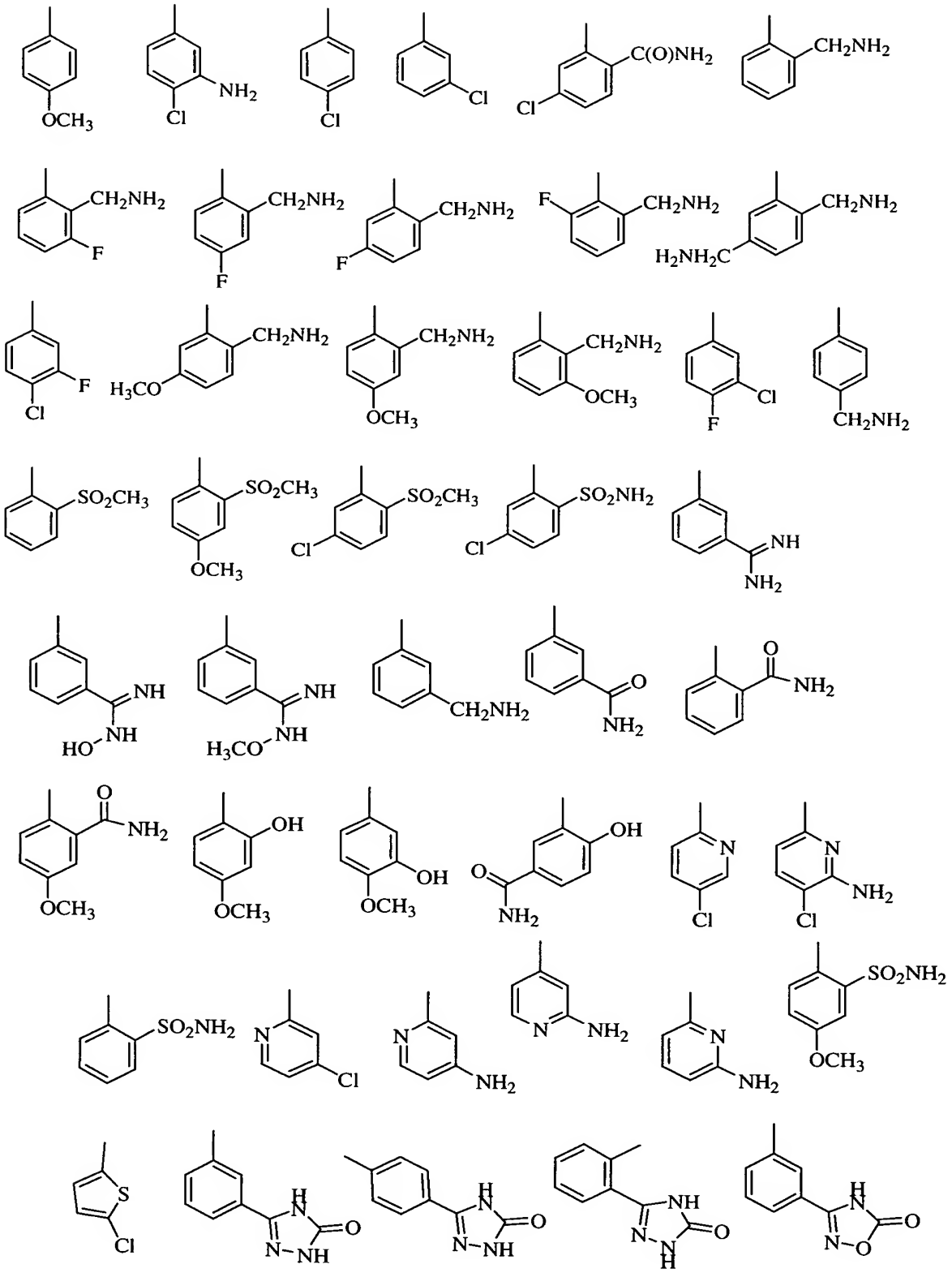
r, at each occurrence, is selected from 0, 1, 2, and 3.

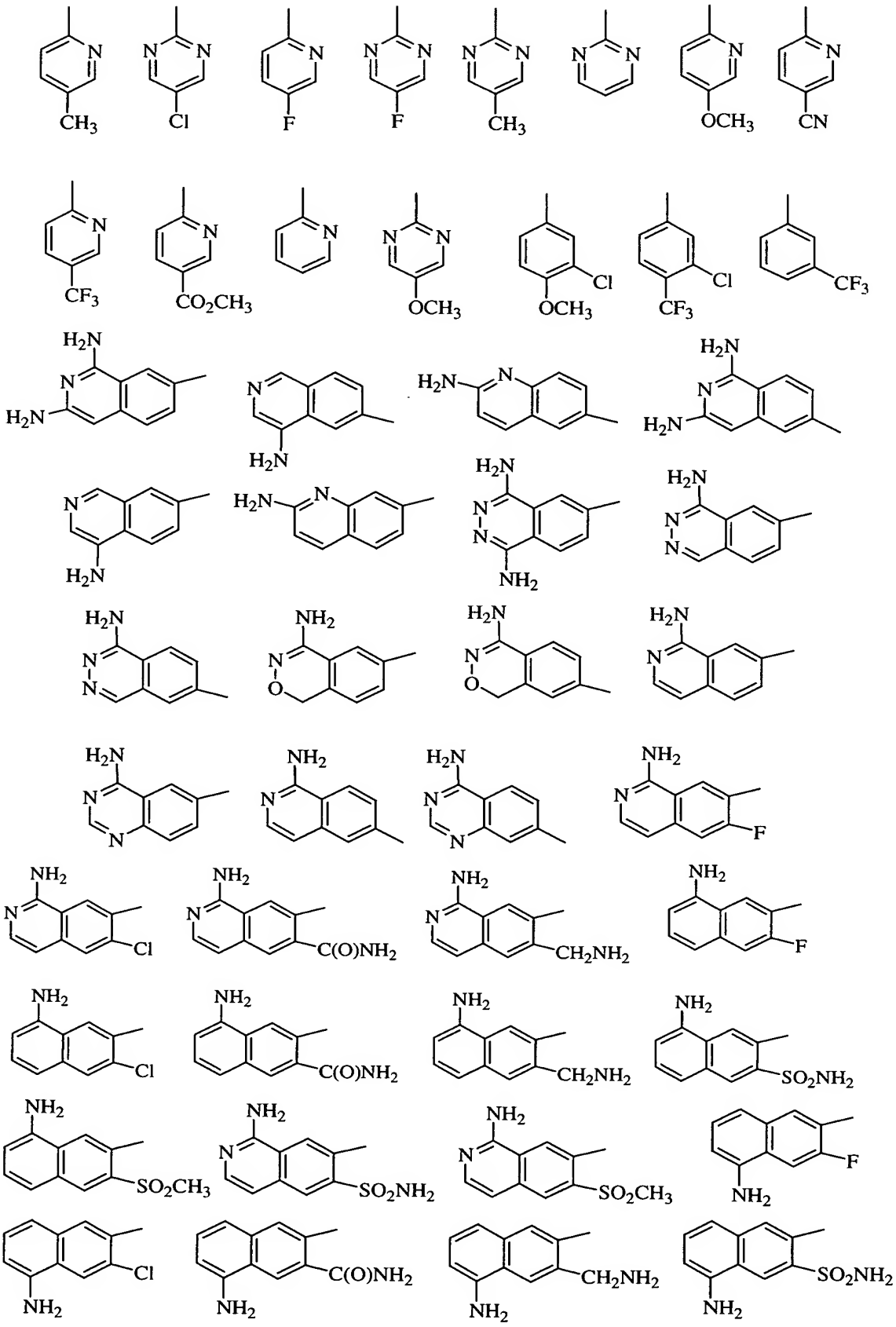
3. A compound according to Claim 2, wherein:

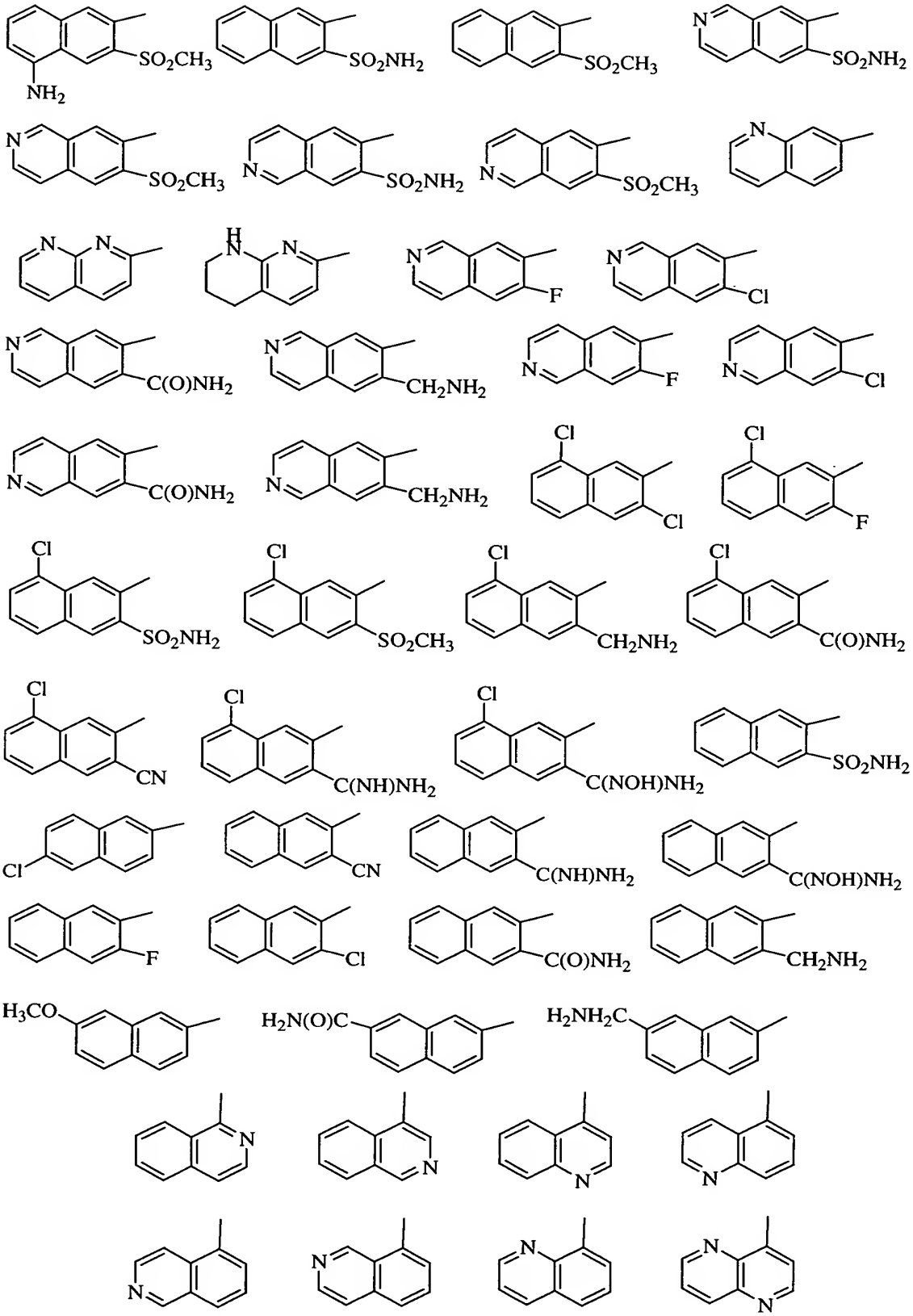
ring M is selected from phenyl, pyrrole, furan, thiophene, pyrazole, imidazole, isoxazole, oxazole, isothiazole, thiazole, pyridine, and pyrimidine;

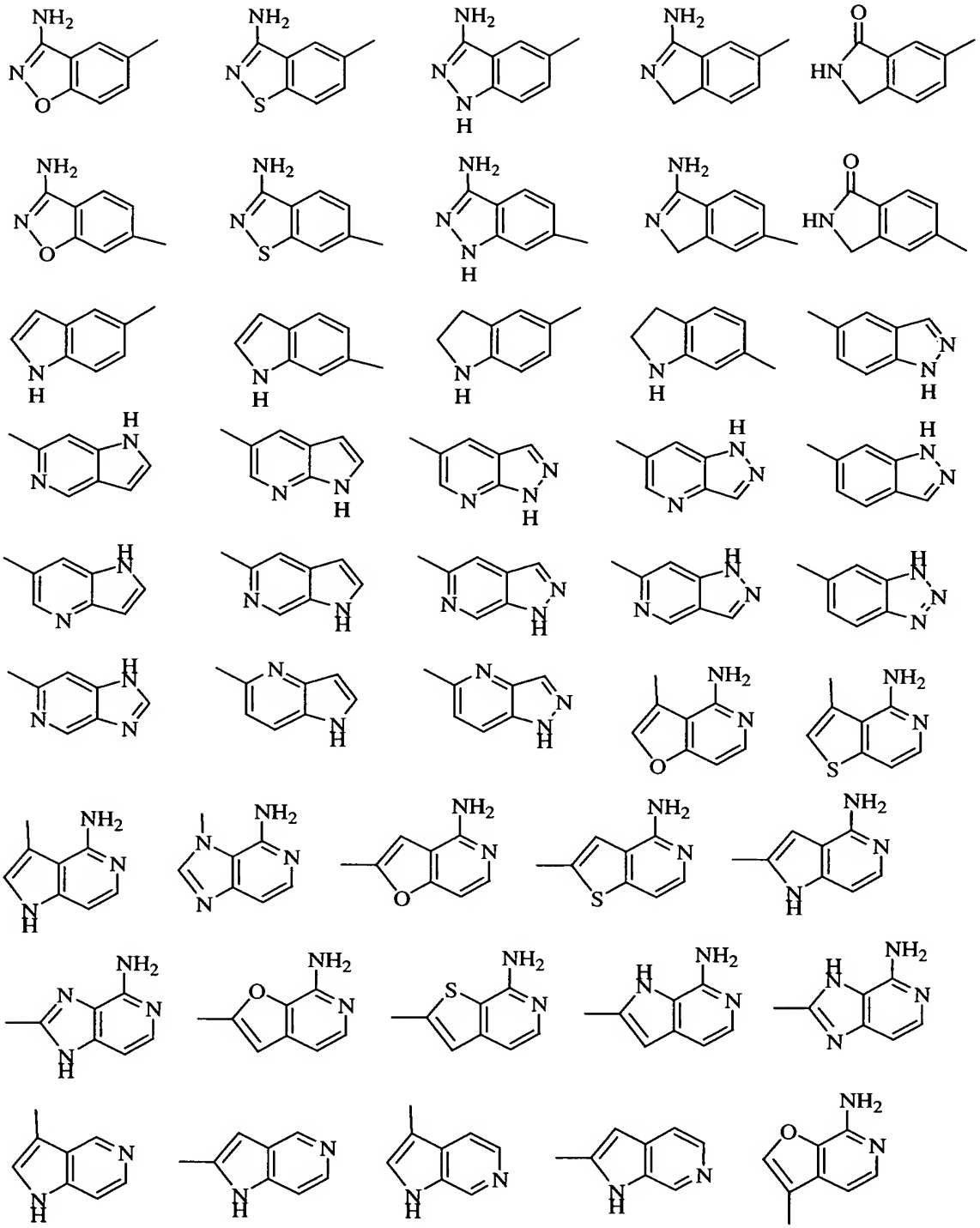
ring M is substituted with 0-2 R^{1a};

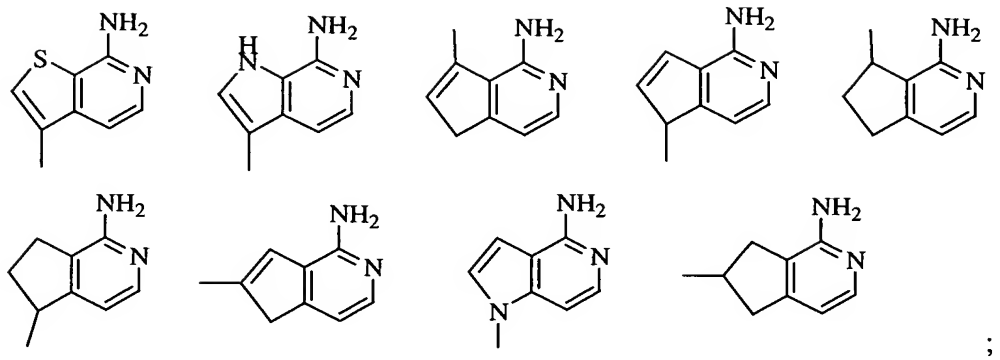
G is selected from the group:











;

R^{1a}, at each occurrence, is selected from H, R^{1b}, $-CH(CH_3)R^{1b}$, $-C(CH_3)_2R^{1b}$, $-CH_2R^{1b}$, and $-CH_2CH_2R^{1b}$, provided that R^{1a} forms other than an N-halo, N-S, or N-CN bond;

alternatively, when two R^{1a} groups are attached to adjacent atoms, together with the atoms to which they are attached, they form a 5-6 membered ring consisting of: carbon atoms and 0-2 heteroatoms selected from the group consisting of N, O, and $S(O)_p$, this ring being substituted with 0-2 R^{4b} and 0-3 ring double bonds;

R^{1b} is selected from H, CH_3, CH_2CH_3, F, Cl, Br, CN, CHO, CF_3, OR^2, $-NR^2R^{2a}$, $-C(O)R^{2b}$, $-CO_2R^{2b}$, $-OC(O)R^2$, $-CO_2R^{2a}$, $-S(O)_pR^2$, $-NR^2(CH_2)_rOR^2$, $-NR^2C(O)R^{2b}$, $-C(O)NR^2R^{2a}$, $-SO_2NR^2R^{2a}$, $-NR^2SO_2R^2$, phenyl substituted with 0-2 R^{4b}, and 5-6 membered aromatic heterocycle consisting of carbon atoms and from 1-4 heteroatoms selected from the group consisting of N, O, and $S(O)_p$ and substituted with 0-2 R^{4b}, provided that R^{1b} forms other than an O-O, N-halo, N-S, or N-CN bond;

R^2, at each occurrence, is selected from H, CF_3, CH_3, CH_2CH_3, $CH_2CH_2CH_3$, $CH(CH_3)_2$, phenyl substituted with 0-2 R^{4b}, a benzyl substituted with 0-2 R^{4b}, C_{3-6} cycloalkyl substituted with 0-2 R^{4b}, C_{3-6} cycloalkyl-CH_2- substituted with 0-2 R^{4b}, and 5-6 membered aromatic heterocycle substituted with 0-2 R^{4b} and consisting of: carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and $S(O)_p$;

R^{2a}, at each occurrence, is selected from H, CF_3, CH_3, CH_2CH_3, $CH_2CH_2CH_3$, $CH(CH_3)_2$, benzyl, phenyl substituted with 0-2 R^{4b}, and 5-6 membered

aromatic heterocycle substituted with 0-2 R^{4b} and consisting of: carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and $S(O)_p$;

R^{2b}, at each occurrence, is selected from CF_3, C_{1-4} alkoxy, CH_3, CH_2CH_3, $CH_2CH_2CH_3$, $CH(CH_3)_2$, benzyl, phenyl substituted with 0-2 R^{4b}, and 5-6 membered aromatic heterocycle substituted with 0-2 R^{4b} and consisting of: carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and $S(O)_p$;

R^{2c}, at each occurrence, is selected from CF_3, OH, OCH_3, OCH_2CH_3, $OCH_2CH_2CH_3$, $OCH(CH_3)_2$, CH_3, CH_2CH_3, $CH_2CH_2CH_3$, $CH(CH_3)_2$, benzyl, phenyl substituted with 0-2 R^{4b}, and 5-6 membered aromatic heterocycle substituted with 0-2 R^{4b} and consisting of carbon atoms and from 1-4 heteroatoms selected from the group consisting of N, O, and $S(O)_p$;

alternatively, R^2 and R^{2a}, together with the nitrogen atom to which they are attached, combine to form a 5 or 6 membered saturated, partially saturated, or unsaturated ring substituted with 0-2 R^{4b} and consisting of: 0-1 additional heteroatoms selected from the group consisting of N, O, and $S(O)_p$;

R^4, at each occurrence, is selected from H, $-(CH_2)_2OR^2$, $-CH_2OR^2$, OR^2, F, Cl, Br, I, CH_3, CH_2CH_3, $CH_2CH_2CH_3$, $CH(CH_3)_2$, $CH_2CH_2CH_2CH_3$, $CH_2CH(CH_3)_2$, $CH(CH_3)CH_2CH_3$, $C(CH_3)_3$, CN, NO_2, $-NR^2R^{2a}$, $-CH_2NR^2R^{2a}$, $-(CH_2)_2NR^2R^{2a}$, $-C(O)R^{2c}$, $-NR^2C(O)R^{2b}$, $-C(O)NR^2R^{2a}$, $-SO_2NR^2R^{2a}$, CF_3, and CF_2CF_3;

R^{4a}, at each occurrence, is selected from H, =O, $-(CH_2)_rOR^2$, $-(CH_2)_r$-F, $-(CH_2)_r$-Br, $-(CH_2)_r$-Cl, C_{1-4} alkyl, $-(CH_2)_r$-CN, $-(CH_2)_rNO_2$, $-(CH_2)_rNR^2R^{2a}$, $-(CH_2)_rC(O)R^{2c}$, $-(CH_2)_rNR^2C(O)R^{2b}$, $-(CH_2)_rC(O)NR^2R^{2a}$, $-(CH_2)_rSO_2NR^2R^{2a}$, $-(CH_2)_rNR^2SO_2NR^2R^{2a}$, $-(CH_2)_rNR^2SO_2$-C_{1-4} alkyl, $-(CH_2)_rC(O)NHSO_2$-C_{1-4} alkyl, $-(CH_2)_rNR^2SO_2R^5$, $-(CH_2)_rS(O)_pR^5$, $-(CH_2)_r(CF_2)_rCF_3$, phenyl substituted with 0-1 R^5, and a 5 membered aromatic heterocycle consisting of: carbon atoms and 1-3 heteroatoms selected from the group consisting of N, O, and $S(O)_p$ substituted with 0-1 R^5;

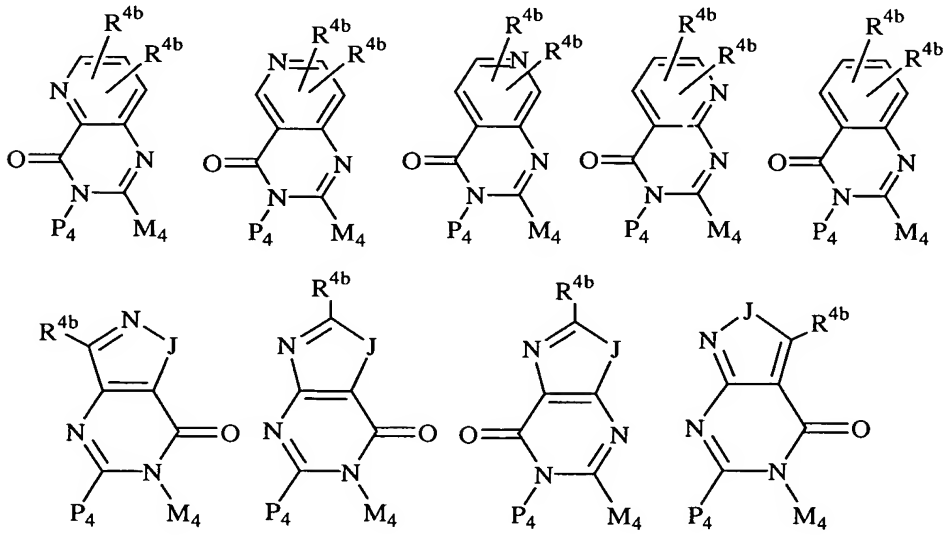
R^{4b}, at each occurrence, is selected from H, =O, OR^3, $-CH_2OR^3$, F, Cl, CH_3, CH_2CH_3, $CH_2CH_2CH_3$, $CH(CH_3)_2$, CN, NO_2, $-NR^3R^{3a}$, $-CH_2NR^3R^{3a}$, $-C(O)R^3$, $-CH_2-C(O)R^3$, $-C(O)OR^{3c}$, $-CH_2-C(O)OR^{3c}$, $-NR^3C(O)R^{3a}$, $-CH_2NR^3C(O)R^{3a}$, $-C(O)NR^3R^{3a}$, $-CH_2-C(O)NR^3R^{3a}$, $-SO_2NR^3R^{3a}$, $-CH_2SO_2NR^3R^{3a}$, $-NR^3SO_2-C_{1-4}$ alkyl, $-CH_2NR^3SO_2-C_{1-4}$ alkyl, $-NR^3SO_2$-phenyl, $-CH_2NR^3SO_2$-phenyl, $-S(O)_pCF_3$, $-CH_2S(O)_pCF_3$, $-S(O)_p-C_{1-4}$ alkyl, $-CH_2S(O)_p-C_{1-4}$ alkyl, $-S(O)_p$-phenyl, $-CH_2S(O)_p$-phenyl, and CF_3;

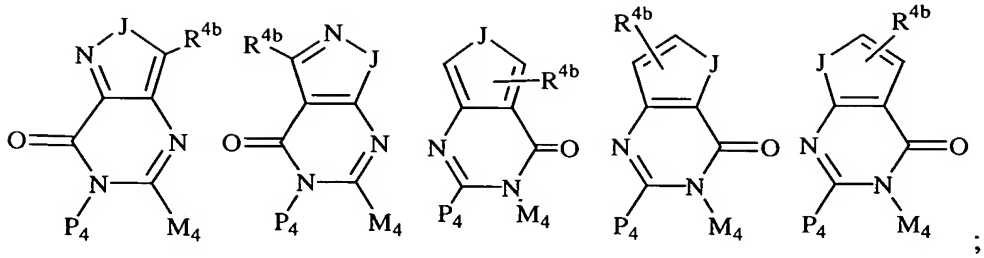
R^5, at each occurrence, is selected from H, =O, CH_3, CH_2CH_3, $CH_2CH_2CH_3$, $CH(CH_3)_2$, OR^3, CH_2OR^3, F, Cl, CN, NO_2, $-NR^3R^{3a}$, $-CH_2NR^3R^{3a}$, $-C(O)R^3$, $-CH_2C(O)R^3$, $-C(O)OR^{3c}$, $-CH_2C(O)OR^{3c}$, $-NR^3C(O)R^{3a}$, $-C(O)NR^3R^{3a}$, $-SO_2NR^3R^{3a}$, $-NR^3SO_2-C_{1-4}$ alkyl, $-NR^3SO_2CF_3$, $-NR^3SO_2$-phenyl, $-S(O)_pCF_3$, $-S(O)_p-C_{1-4}$ alkyl, $-S(O)_p$-phenyl, CF_3, phenyl substituted with 0-2 R^6, naphthyl substituted with 0-2 R^6, and benzyl substituted with 0-2 R^6;

R^6, at each occurrence, is selected from H, OH, OR^2, F, Cl, CH_3, CH_2CH_3, $CH_2CH_2CH_3$, $CH(CH_3)_2$, CN, NO_2, $-NR^2R^{2a}$, $-CH_2NR^2R^{2a}$, $-C(O)R^{2b}$, $-CH_2C(O)R^{2b}$, $-NR^2C(O)R^{2b}$, $-SO_2NR^2R^{2a}$, and $NR^2SO_2C_{1-4}$ alkyl; and

r, at each occurrence, is selected from 0, 1, and 2.

4. A compound according to Claim 3, wherein the compound is selected from:

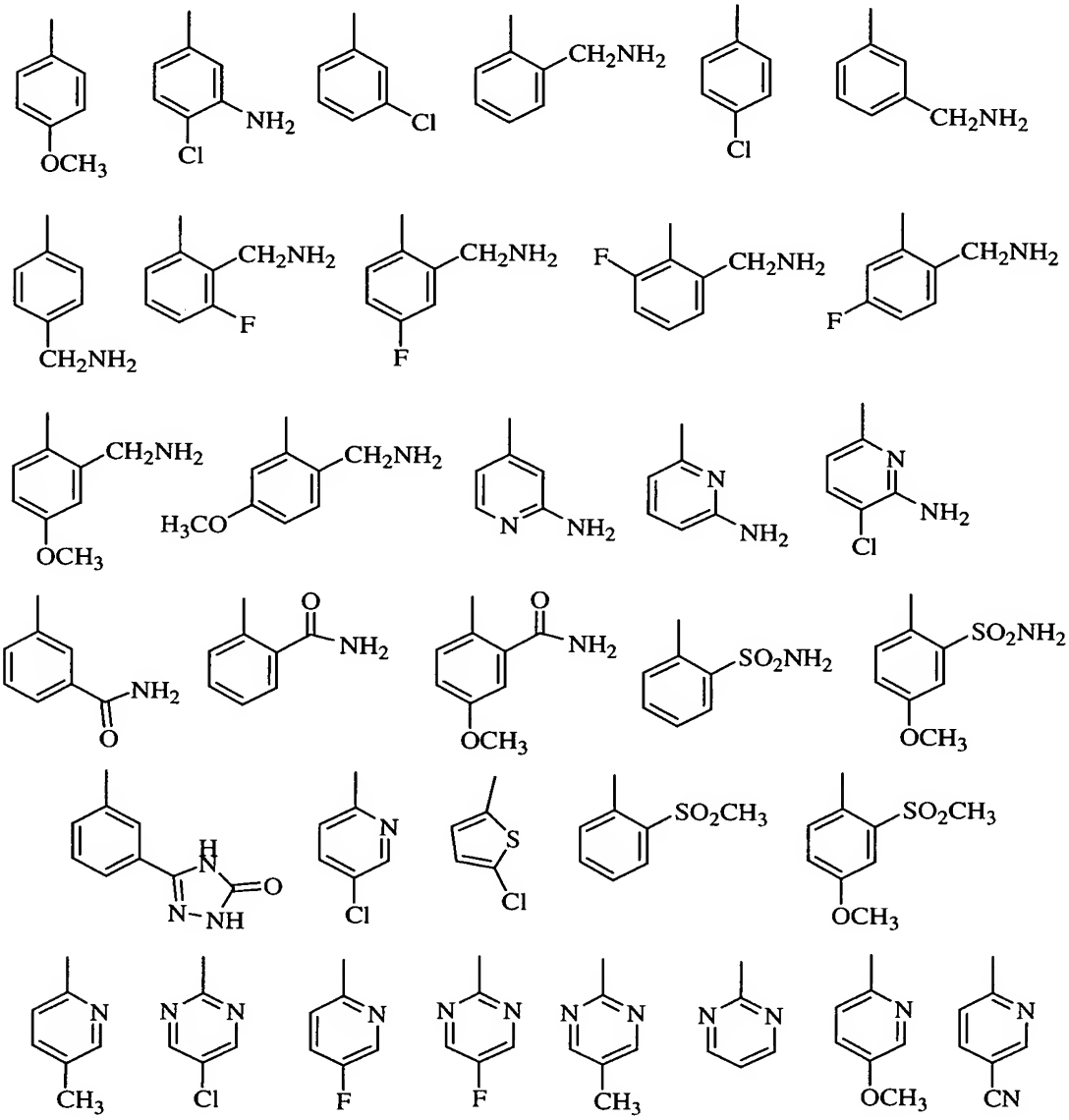


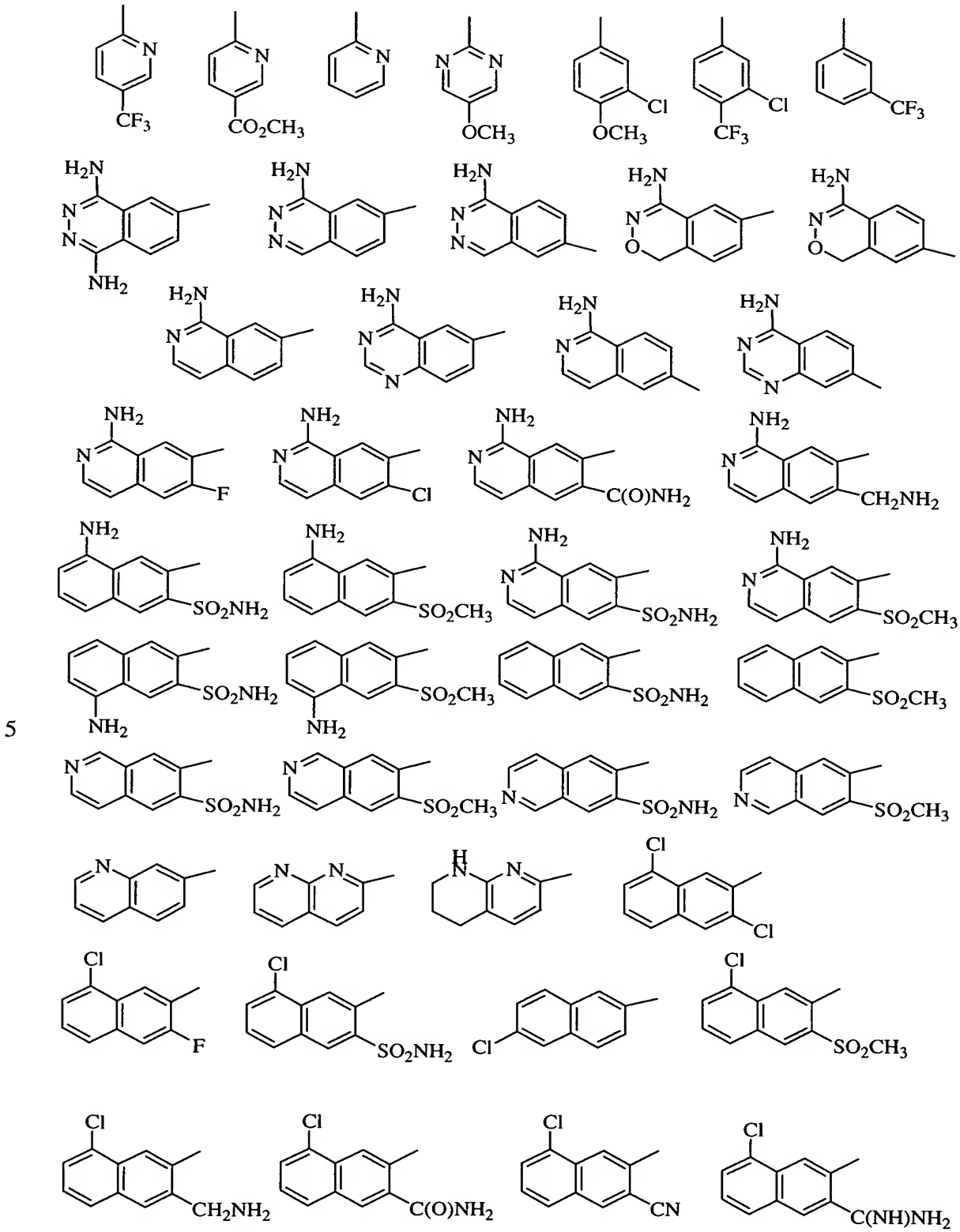


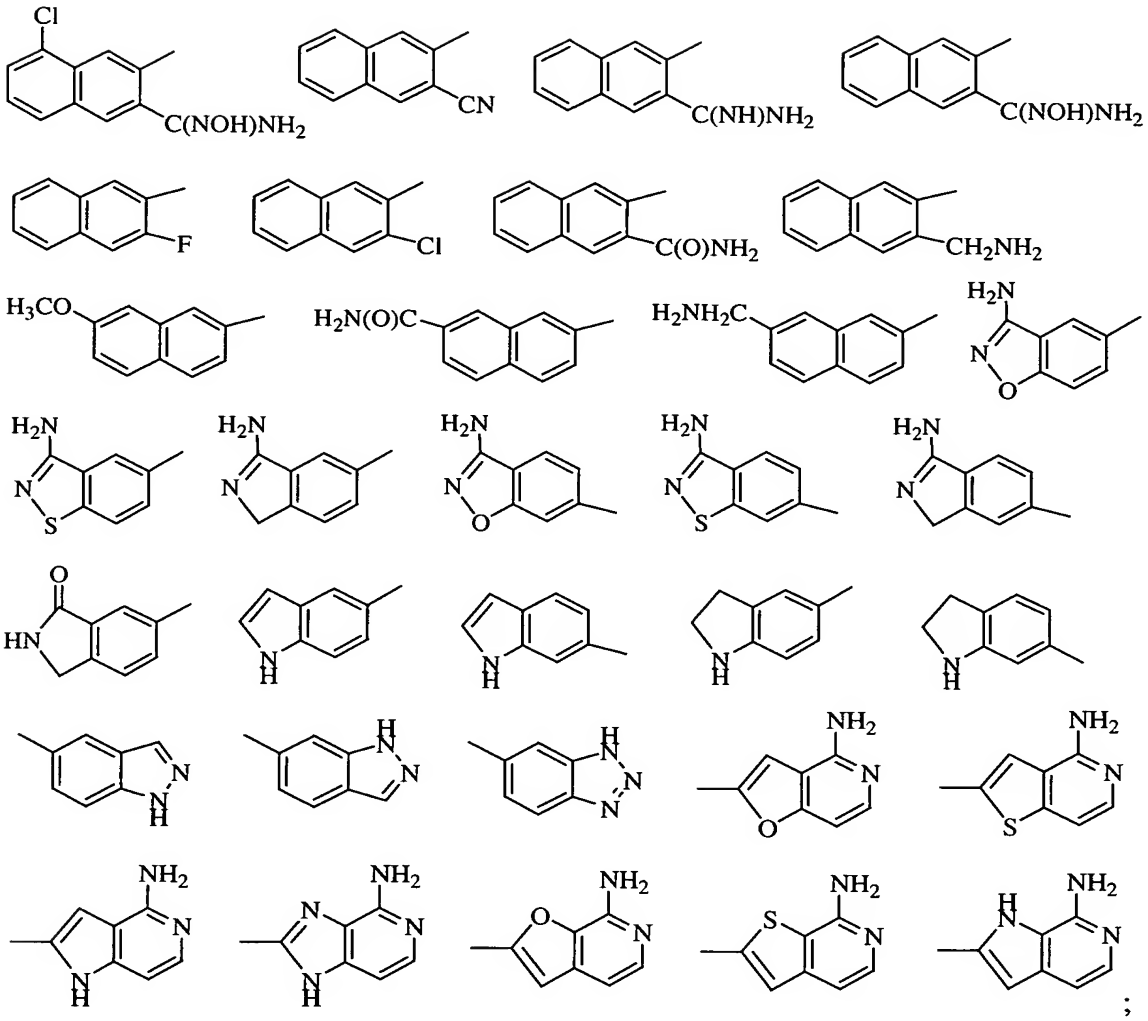
;

J is selected from O, S, NH, and NR^{1a};

G is selected from the group:







;

R^{1a}, at each occurrence, is selected from H, R^{1b}, $-CH(CH_3)R^{1b}$, $-C(CH_3)_2R^{1b}$, and $-CH_2R^{1b}$, provided that R^{1a} forms other than an N-halo, N-S, or N-CN bond;

R^{1b} is selected from CH_3, CH_2CH_3, F, Cl, Br, -CN, CF_3, OR^2, $-NR^2R^{2a}$, $-C(O)R^{2b}$, $-CO_2R^{2b}$, $-CO_2R^{2a}$, $-S(O)_pR^2$, $-C(O)NR^2R^{2a}$, $-SO_2NR^2R^{2a}$, $-NR^2SO_2R^2$, and 5-6 membered aromatic heterocycle consisting of carbon atoms and from 1-4 heteroatoms selected from the group consisting of N, O, and $S(O)_p$ and substituted with 0-2 R^{4b}, provided that R^{1b} forms other than an O-O, N-halo, N-S, or N-CN bond;

R^2, at each occurrence, is selected from H, CH_3, CH_2CH_3, $CH_2CH_2CH_3$, $CH(CH_3)_2$, phenyl substituted with 0-1 R^{4b}, benzyl substituted with 0-1 R^{4b},

C_{3-5} cycloalkyl substituted with 0-1 R^{4b}, C_{3-5} cycloalkyl-CH_2- substituted with 0-1 R^{4b}, and 5-6 membered aromatic heterocycle substituted with 0-1 R^{4b} and consisting of: carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and $S(O)_p$;

R^{2a}, at each occurrence, is selected from H, CH_3, CH_2CH_3, $CH_2CH_2CH_3$, $CH(CH_3)_2$, benzyl, phenyl substituted with 0-1 R^{4b}, and 5-6 membered aromatic heterocycle substituted with 0-1 R^{4b} and consisting of: carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and $S(O)_p$;

alternatively, R^2 and R^{2a}, together with the nitrogen atom to which they are attached, combine to form a 5 or 6 membered saturated, partially saturated, or unsaturated ring substituted with 0-1 R^{4b} and consisting of: 0-1 additional heteroatoms selected from the group consisting of N, O, and $S(O)_p$;

R^{2b}, at each occurrence, is selected from OCH_3, $-OCH_2CH_3$, $OCH_2CH_2CH_3$, $OCH(CH_3)_2$, CH_3, CH_2CH_3, $CH_2CH_2CH_3$, $CH(CH_3)_2$, benzyl, phenyl substituted with 0-1 R^{4b}, and 5-6 membered aromatic heterocycle substituted with 0-1 R^{4b} and consisting of: carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and $S(O)_p$;

R^{2c}, at each occurrence, is selected from OH, OCH_3, OCH_2CH_3, $OCH_2CH_2CH_3$, $OCH(CH_3)_2$, CH_3, CH_2CH_3, $CH_2CH_2CH_3$, $CH(CH_3)_2$, benzyl, phenyl substituted with 0-1 R^{4b}, and 5-6 membered aromatic heterocycle substituted with 0-1 R^{4b} and consisting of carbon atoms and from 1-4 heteroatoms selected from the group consisting of N, O, and $S(O)_p$;

R^4, at each occurrence, is selected from OH, OR^2, CH_2OR^2, $(CH_2)_2OR^2$, F, Br, Cl, I, CH_3, CH_2CH_3, $CH_2CH_2CH_3$, $CH(CH_3)_2$, $CH_2CH_2CH_2CH_3$, $CH_2CH(CH_3)_2$, $CH(CH_3)CH_2CH_3$, $C(CH_3)_3$, $-NR^2R^{2a}$, $-CH_2NR^2R^{2a}$, $-(CH_2)_2NR^2R^{2a}$, CF_3, and CF_2CF_3;

R^{4a}, at each occurrence, is selected from H, =O, $-(CH_2)_rOR^2$, F, Br, Cl, C_{1-4} alkyl, $-(CH_2)_rNR^2R^{2a}$, $-(CH_2)_rC(O)R^{2c}$, $-(CH_2)_rNR^2C(O)R^{2b}$, $-(CH_2)_rC(O)NR^2R^{2a}$, $-(CH_2)_rSO_2NR^2R^{2a}$, $-(CH_2)_rNR^2SO_2R^5$, $-(CH_2)_rS(O)_pR^5$,

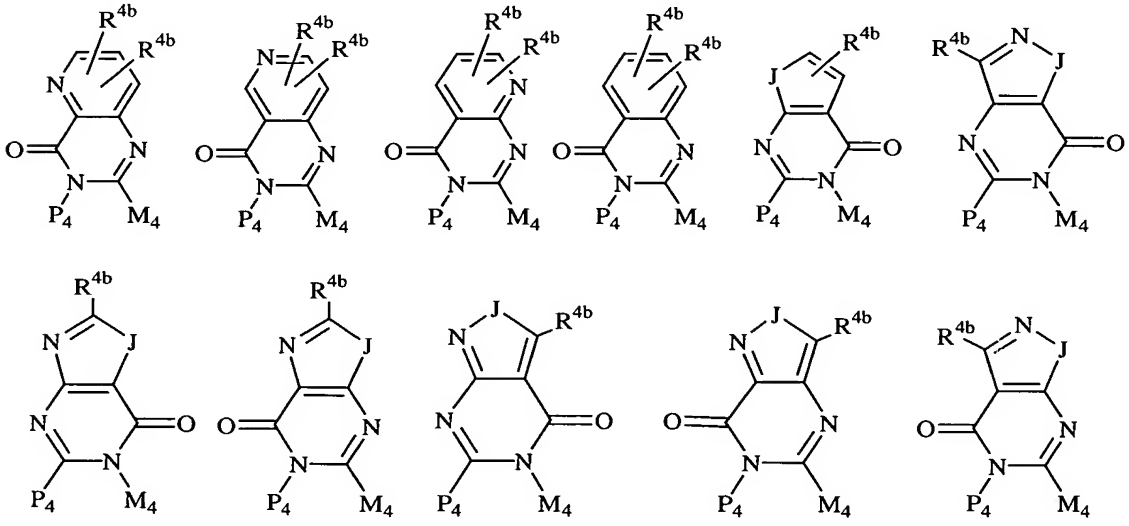
$-(CH_2)_r(CF_2)_rCF_3$, phenyl substituted with 0-1 R^5, and a 5 membered aromatic heterocycle consisting of: carbon atoms and 1-3 N and is substituted with 1 R^5;

R^{4b}, at each occurrence, is selected from H, =O, OR^3, $-CH_2OR^3$, F, Cl, CH_3, CH_2CH_3, $CH_2CH_2CH_3$, $CH(CH_3)_2$, CN, NO_2, $-NR^3R^{3a}$, $-CH_2NR^3R^{3a}$, $-C(O)R^3$, $-C(O)OR^{3c}$, $-NR^3C(O)R^{3a}$, $-C(O)NR^3R^{3a}$, $-SO_2NR^3R^{3a}$, $-NR^3SO_2-C_{1-4}$ alkyl, $-NR^3SO_2$-phenyl, $-S(O)_p-C_{1-4}$ alkyl, $-S(O)_p$-phenyl, and CF_3;

R^5, at each occurrence, is selected from H, =O, CH_3, CH_2CH_3, $CH_2CH_2CH_3$, $CH(CH_3)_2$, OR^3, CH_2OR^3, F, Cl, CN, NO_2, $-NR^3R^{3a}$, $-CH_2NR^3R^{3a}$, $-C(O)R^3$, $-C(O)OR^{3c}$, $-NR^3C(O)R^{3a}$, $-C(O)NR^3R^{3a}$, $-SO_2NR^3R^{3a}$, $-NR^3SO_2-C_{1-4}$ alkyl, $-NR^3SO_2$-phenyl, $-S(O)_p-C_{1-4}$ alkyl, $-S(O)_p$-phenyl, CF_3, phenyl substituted with 0-2 R^6, naphthyl substituted with 0-2 R^6, and benzyl substituted with 0-2 R^6; and

R^6, at each occurrence, is selected from H, OH, OR^2, F, Cl, CH_3, CH_2CH_3, $CH_2CH_2CH_3$, $CH(CH_3)_2$, CN, NO_2, $-NR^2R^{2a}$, $-CH_2NR^2R^{2a}$, $-C(O)R^{2b}$, $-CH_2C(O)R^{2b}$, $-NR^2C(O)R^{2b}$, and $-SO_2NR^2R^{2a}$.

5. A compound according to Claim 4, wherein the compound is selected from:

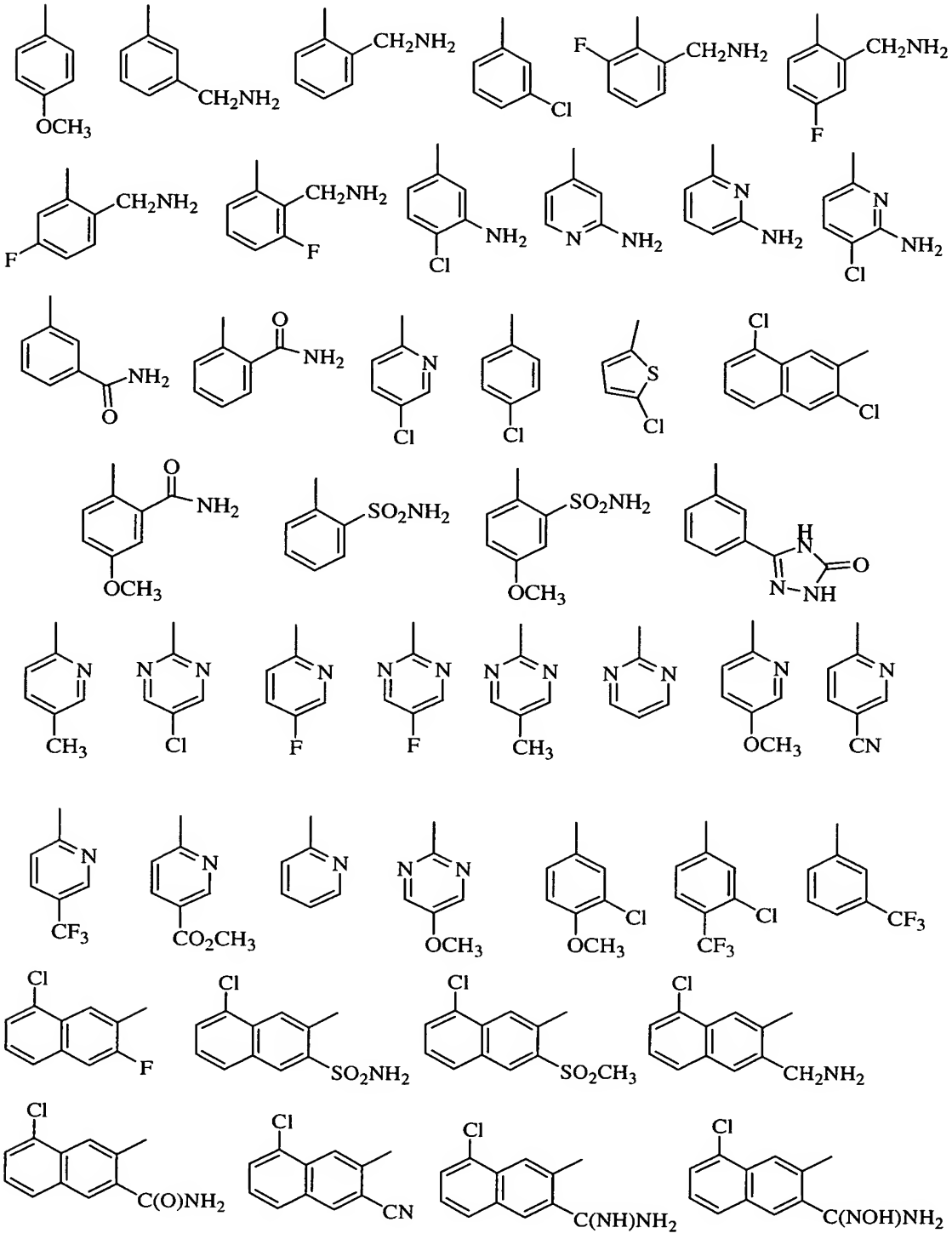


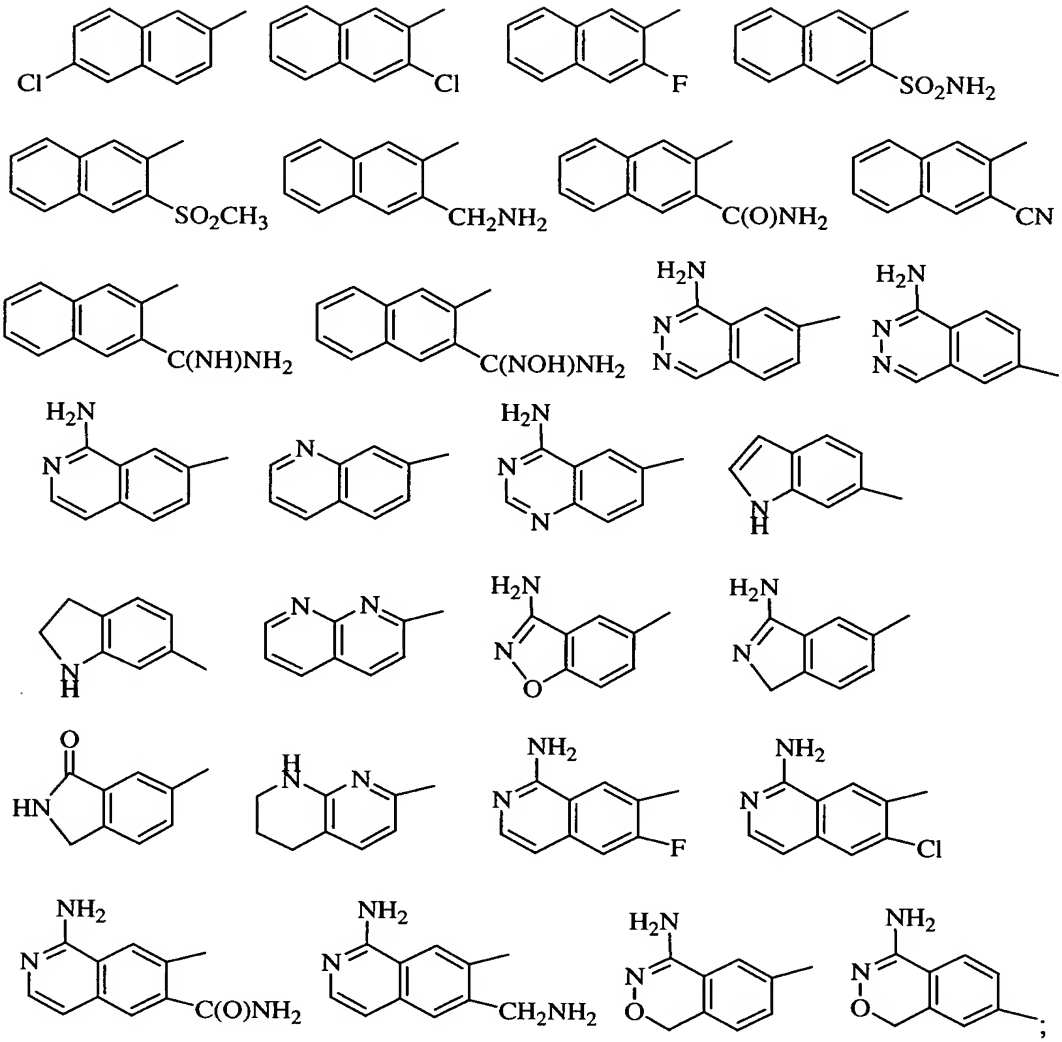
J is selected from O, S, NH, and NR^{1a};

P_4 is $-G_1$-G;

M_4 is –Z-A-B;

G is selected from:





G_1 is absent or is selected from CH_2CH_2, CH_2O, OCH_2, CH_2NH, $NHCH_2$, $CH_2C(O)$, $C(O)CH_2$, $C(O)NH$, $NHC(O)$, SO_2NH, and $NHSO_2$, provided that G_1 does not form a N-S, NCH_2N, NCH_2O, or NCH_2S bond with either group to which it is attached;

A is selected from phenyl, pyridyl, and pyrimidyl, and is substituted with 0-2 R^4;

B is selected from phenyl, pyrrolidinyl, N-pyrrolidino-carbonyl, morpholinyl, N-morpholino-carbonyl, 1,2,3-triazolyl, imidazolyl, and benzimidazolyl, and is substituted with 0-1 R^{4a};

R^{1a}, at each occurrence, is selected from H, CH_3, CH_2CH_3, $CH_2CH_2CH_3$, CH_2F, CH_2Cl, Br, CH_2Br, -CN, CH_2CN, CF_3, CH_2CF_3, OCH_3, CH_2OH,

$C(CH_3)_2OH$, CH_2OCH_3, NH_2, CH_2NH_2, $NHCH_3$, CH_2NHCH_3, $N(CH_3)_2$, $CH_2N(CH_3)_2$, CO_2H, $COCH_3$, CO_2CH_3, $CH_2CO_2CH_3$, SCH_3, CH_2SCH_3, $S(O)CH_3$, $CH_2S(O)CH_3$, $S(O)_2CH_3$, $CH_2S(O)_2CH_3$, $C(O)NH_2$, $CH_2C(O)NH_2$, SO_2NH_2, $CH_2SO_2NH_2$, $NHSO_2CH_3$, $CH_2NHSO_2CH_3$, pyridin-2-yl, pyridin-3-yl, pyridin-4-yl, pyridin-2-yl-N-oxide, pyridin-3-yl-N-oxide, pyridin-4-yl-N-oxide, imidazol-1-yl, CH_2-imidazol-1-yl, 4-methyl-oxazol-2-yl, 4-N,N-dimethylaminomethyl-oxazol-2-yl, 1,2,3,4-tetrazol-1-yl, 1,2,3,4-tetrazol-5-yl, -CH_2-1,2,3,4-tetrazol-1-yl, and -CH_2-1,2,3,4-tetrazol-5-yl, provided that R^{1a} forms other than an N-halo, N-S, or N-CN bond;

R^2, at each occurrence, is selected from H, CH_3, CH_2CH_3, cyclopropylmethyl, cyclobutyl, and cyclopentyl;

R^{2a}, at each occurrence, is H or CH_3;

alternatively, R^2 and R^{2a}, together with the atom to which they are attached, combine to form pyrrolidine substituted with 0-2 R^{4b} or piperidine substituted with 0-2 R^{4b};

R^{2b}, at each occurrence, is selected from OCH_3, OCH_2CH_3, CH_3, and CH_2CH_3;

R^{2c}, at each occurrence, is selected from OH, OCH_3, OCH_2CH_3, CH_3, and CH_2CH_3;

R^4, at each occurrence, is selected from OH, OR^2, CH_2OR^2, $(CH_2)_2OR^2$, F, Br, Cl, I, CH_3, CH_2CH_3, $CH_2CH_2CH_3$, $CH(CH_3)_2$, $CH_2CH_2CH_2CH_3$, $CH_2CH(CH_3)_2$, $CH(CH_3)CH_2CH_3$, $C(CH_3)_3$, -NR^2R^{2a}, -$CH_2NR^2R^{2a}$, -$(CH_2)_2NR^2R^{2a}$, CF_3, and CF_2CF_3;

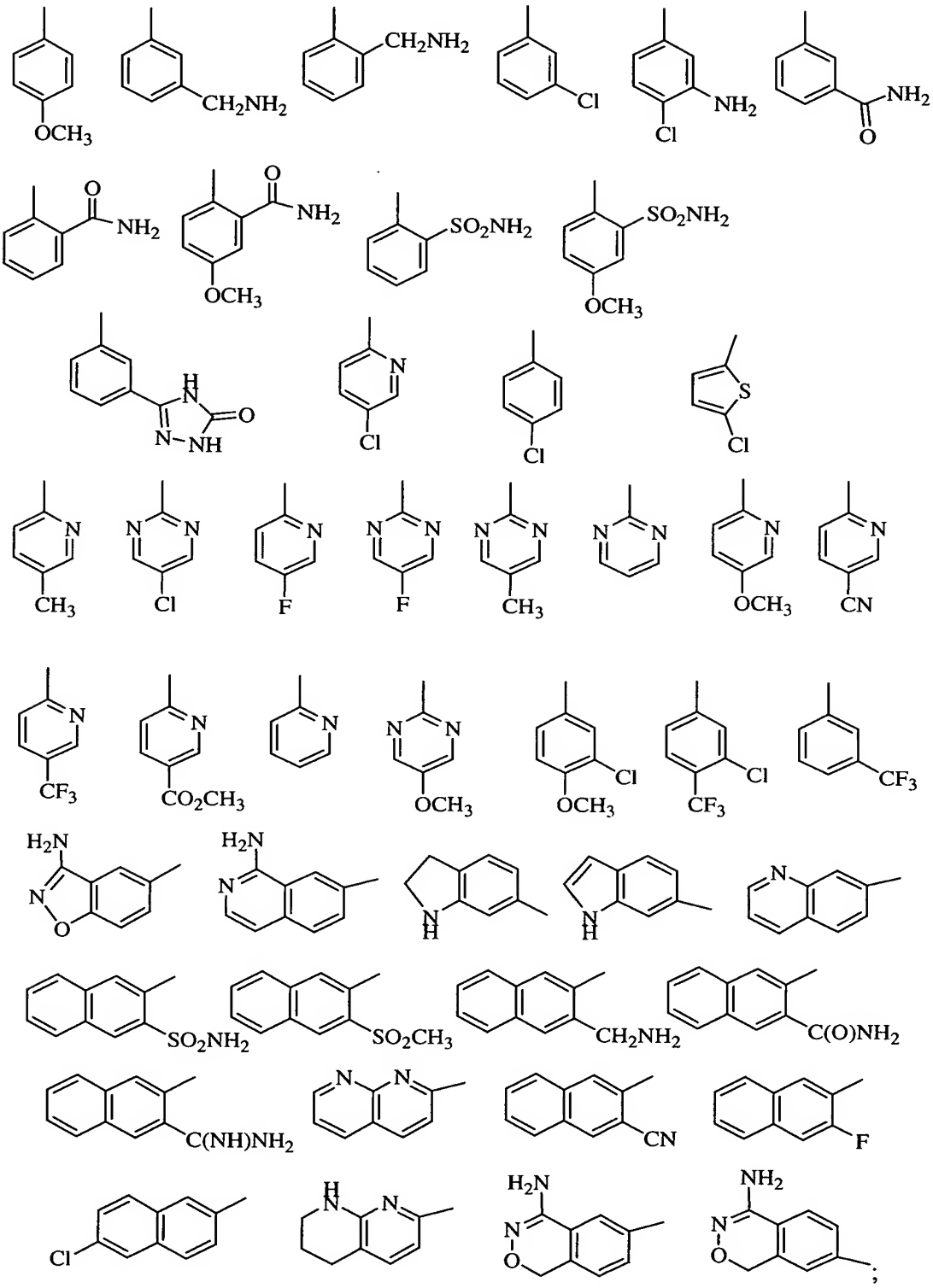
R^{4a} is selected from $C_{1\text{-}4}$ alkyl, CF_3, OR^2, -CH_2OR^2, -$(CH_2)_2OR^2$, -NR^2R^{2a}, -$CH_2NR^2R^{2a}$, -$(CH_2)_2NR^2R^{2a}$, -$S(O)_pR^5$, -$SO_2NR^2R^{2a}$, and 1-CF_3-tetrazol-2-yl;

R^{4b}, at each occurrence, is selected from H, CH_3, and OH; and

R^5, at each occurrence, is selected from CF_3, CH_3, CH_2CH_3, $CH_2CH_2CH_3$, $CH(CH_3)_2$, phenyl, and benzyl.

6. A compound according to Claim 5, wherein the compound is selected from:

G is selected from:



A is selected from the group: phenyl, 2-pyridyl, 3-pyridyl, 2-pyrimidyl, 2-Cl-phenyl, 3-Cl-phenyl, 2-F-phenyl, 3-F-phenyl, 2-methylphenyl, 2-aminophenyl, and 2-methoxyphenyl; and

B is selected from the group: 2-(aminosulfonyl)phenyl, 2-(methylaminosulfonyl)phenyl, N-pyrrolidino-carbonyl, 2-(methylsulfonyl)phenyl, 2-(N,N-dimethylaminomethyl)phenyl, 2-(N-methylaminomethyl)phenyl, 2-(N-ethyl-N-methylaminomethyl)phenyl, 2-(N-pyrrolidinylmethyl)phenyl, 1-methyl-2-imidazolyl, 2-methyl-1-imidazolyl, 2-(dimethylaminomethyl)-1-imidazolyl, 2-(methylaminomethyl)-1-imidazolyl, 2-(N-(cyclopropylmethyl)aminomethyl)phenyl, 2-(N-(cyclobutyl)aminomethyl)phenyl, 2-(N-(cyclopentyl)aminomethyl)phenyl, 2-(N-(4-hydroxypiperidinyl)methyl)phenyl, and 2-(N-(3-hydroxypyrrolidinyl)methyl)phenyl.

7. A compound according to Claim 6, wherein the compound is selected from the group:

2-biphenyl-4-yl-6-chloro-3-(4-chloro-phenyl)-3*H*-quinazolin-4-one;

6-chloro-3-(4-chloro-phenyl)-2-phenyl-3*H*-quinazolin-4-one;

3-(5-chloro-pyridin-2-yl)-2-[4-(1H-pyrrol-2-yl)-phenyl]-3H-quinazolin-4-one;

3-(5-chloro-pyridin-2-yl)-2-[4-(1-methyl-1H-pyrrol-2-yl)-phenyl]-3H-quinazolin-4-one;

3-(5-chloro-pyridin-2-yl)-2-[4-(1-ethyl-1H-pyrrol-2-yl)-phenyl]-3H-quinazolin-4-one;

2-{4-[1-(2-amino-ethyl)-1H-pyrrol-2-yl]-phenyl}-3-(5-chloro-pyridin-2-yl)-3H-quinazolin-4-one;

3-(5-chloro-pyridin-2-yl)-2-{4-[1-(2-methylamino-ethyl)-1H-pyrrol-2-yl]-phenyl}-3H-quinazolin-4-one;

3-(5-chloro-pyridin-2-yl)-2-{4-[1-(2-ethylamino-ethyl)-1H-pyrrol-2-yl]-phenyl}-3H-quinazolin-4-one;

2-{4-[1-(2-benzylamino-ethyl)-1H-pyrrol-2-yl]-phenyl}-3-(5-chloro-pyridin-2-yl)-3H-quinazolin-4-one;

3-pyridin-2-yl-2-[4-(1-{2-[(pyridin-2-ylmethyl)-amino]-ethyl}-1H-pyrrol-2-yl)-phenyl]-3H-quinazolin-4-one;

3-pyridin-2-yl-2-[4-(1-{2-[(pyridin-3-ylmethyl)-amino]-ethyl}-1H-pyrrol-2-yl)-phenyl]-3H-quinazolin-4-one;

3-pyridin-2-yl-2-[4-(1-{2-[(pyridin-4-ylmethyl)-amino]-ethyl}-1H-pyrrol-2-yl)-phenyl]-3H-quinazolin-4-one;

2-[4-(1-benzyl-1H-pyrrol-2-yl)-phenyl]-3-(5-chloro-pyridin-2-yl)-3H-quinazolin-4-one;

3-(5-chloro-pyridin-2-yl)-2-[4-(1-pyridin-2-ylmethyl-1H-pyrrol-2-yl)-phenyl]-3H-quinazolin-4-one;

3-(5-chloro-pyridin-2-yl)-2-[4-(1-pyridin-3-ylmethyl-1H-pyrrol-2-yl)-phenyl]-3H-quinazolin-4-one;

3-(5-chloro-pyridin-2-yl)-2-[4-(1-pyridin-4-ylmethyl-1H-pyrrol-2-yl)-phenyl]-3H-quinazolin-4-one;

3-(5-chloro-pyridin-2-yl)-2-[4-(1-cyclohexyl-1H-pyrrol-2-yl)-phenyl]-3H-quinazolin-4-one;

3-(5-chloro-pyridin-2-yl)-2-{4-[1-(tetrahydro-pyran-4-yl)-1H-pyrrol-2-yl]-phenyl}-3H-quinazolin-4-one;

3-(5-chloro-pyridin-2-yl)-2-[4-(1-piperidin-4-yl-1H-pyrrol-2-yl)-phenyl]-3H-quinazolin-4-one;

3-(5-chloro-pyridin-2-yl)-2-{4-[1-(1-methyl-piperidin-4-yl)-1H-pyrrol-2-yl]-phenyl}-3H-quinazolin-4-one;

2-{4-[1-(1-acetyl-piperidin-4-yl)-1H-pyrrol-2-yl]-phenyl}-3-(5-chloro-pyridin-2-yl)-3H-quinazolin-4-one;

3-(5-chloro-pyridin-2-yl)-2-[4-(1-isopropyl-1H-pyrrol-2-yl)-phenyl]-3H-quinazolin-4-one;

3-(5-chloro-pyridin-2-yl)-2-[4-(1-cyclopropyl-1H-pyrrol-2-yl)-phenyl]-3H-quinazolin-4-one;

3-(5-chloro-pyridin-2-yl)-2-[4-(1-cyclobutyl-1H-pyrrol-2-yl)-phenyl]-3H-quinazolin-4-one;

3-(5-chloro-pyridin-2-yl)-2-[4-(1-cyclopentyl-1H-pyrrol-2-yl)-phenyl]-3H-quinazolin-4-one;

3-(5-chloro-pyridin-2-yl)-2-{4-[1-(tetrahydro-furan-3-yl)-1H-pyrrol-2-yl]-phenyl}-3H-quinazolin-4-one;

3-(5-chloro-pyridin-2-yl)-2-[4-(2',3',4',5'-tetrahydro-1'H-[1,3']bipyrrolyl-2-yl)-phenyl]-3H-quinazolin-4-one;

3-(5-chloro-pyridin-2-yl)-2-[4-(1'-methyl-2',3',4',5'-tetrahydro-1'H-[1,3']bipyrrolyl-2-yl)-phenyl]-3H-quinazolin-4-one;

2-[4-(1'-acetyl-2',3',4',5'-tetrahydro-1'H-[1,3']bipyrrolyl-2-yl)-phenyl]-3-(5-chloro-pyridin-2-yl)-3H-quinazolin-4-one;

3-(5-chloro-pyridin-2-yl)-2-[4-(1'-methanesulfonyl-2',3',4',5'-tetrahydro-1'H-[1,3']bipyrrolyl-2-yl)-phenyl]-3H-quinazolin-4-one;

N-[2-(2-{4-[3-(5-chloro-pyridin-2-yl)-4-oxo-3,4-dihydro-quinazolin-2-yl]-phenyl}-pyrrol-1-yl)-ethyl]-acetamide;

3-(5-chloro-pyridin-2-yl)-2-{4-[1-(2-hydroxy-ethyl)-1H-pyrrol-2-yl]-phenyl}-3H-quinazolin-4-one;

3-(5-chloro-pyridin-2-yl)-2-{4-[1-(2-methoxy-ethyl)-1H-pyrrol-2-yl]-phenyl}-3H-quinazolin-4-one;

3-(5-chloro-pyridin-2-yl)-2-{4-[1-(2-methoxy-1-methyl-ethyl)-1H-pyrrol-2-yl]-phenyl}-3H-quinazolin-4-one;

2-(2-{4-[3-(5-chloro-pyridin-2-yl)-4-oxo-3,4-dihydro-quinazolin-2-yl]-phenyl}-pyrrol-1-yl)-acetamide;

2-(2-{4-[3-(5-chloro-pyridin-2-yl)-4-oxo-3,4-dihydro-quinazolin-2-yl]-phenyl}-pyrrol-1-yl)-N-methyl-acetamide;

3-(5-chloro-pyridin-2-yl)-2-[4-(1H-imidazol-2-yl)-phenyl]-3H-quinazolin-4-one;

3-(5-chloro-pyridin-2-yl)-2-[4-(1-methyl-1H-imidazol-2-yl)-phenyl]-3H-quinazolin-4-one;

3-(5-chloro-pyridin-2-yl)-2-[4-(1-ethyl-1H-imidazol-2-yl)-phenyl]-3H-quinazolin-4-one;

2-{4-[1-(2-amino-ethyl)-1H-imidazol-2-yl]-phenyl}-3-(5-chloro-pyridin-2-yl)-3H-quinazolin-4-one;

3-(5-chloro-pyridin-2-yl)-2-{4-[1-(2-methylamino-ethyl)-1H-imidazol-2-yl]-phenyl}-3H-quinazolin-4-one;

2-{4-[1-(2-ethylamino-ethyl)-1H-imidazol-2-yl]-phenyl}-3-pyridin-2-yl-3H-quinazolin-4-one;

2-{4-[1-(2-benzylamino-ethyl)-1H-imidazol-2-yl]-phenyl}-3-pyridin-2-yl-3H-quinazolin-4-one;

3-pyridin-2-yl-2-[4-(1-{2-[(pyridin-2-ylmethyl)-amino]-ethyl}-1H-imidazol-2-yl)-phenyl]-3H-quinazolin-4-one;

3-pyridin-2-yl-2-[4-(1-{2-[(pyridin-3-ylmethyl)-amino]-ethyl}-1H-imidazol-2-yl)-phenyl]-3H-quinazolin-4-one;

2-{4-[1-(2-benzylamino-ethyl)-1H-imidazol-2-yl]-phenyl}-3-pyridin-2-yl-3H-quinazolin-4-one;

3-(5-chloro-pyridin-2-yl)-2-[4-(1-pyridin-2-ylmethyl-1H-imidazol-2-yl)-phenyl]-3H-quinazolin-4-one;

3-(5-chloro-pyridin-2-yl)-2-[4-(1-pyridin-3-ylmethyl-1H-imidazol-2-yl)-phenyl]-3H-quinazolin-4-one;

3-(5-chloro-pyridin-2-yl)-2-[4-(1-pyridin-4-ylmethyl-1H-imidazol-2-yl)-phenyl]-3H-quinazolin-4-one;

3-(5-chloro-pyridin-2-yl)-2-[4-(1-cyclohexyl-1H-imidazol-2-yl)-phenyl]-3H-quinazolin-4-one;

3-(5-chloro-pyridin-2-yl)-2-{4-[1-(tetrahydro-pyran-4-yl)-1H-imidazol-2-yl]-phenyl}-3H-quinazolin-4-one;

3-(5-chloro-pyridin-2-yl)-2-[4-(1-piperidin-4-yl-1H-imidazol-2-yl)-phenyl]-3H-quinazolin-4-one;

3-(5-chloro-pyridin-2-yl)-2-{4-[1-(1-methyl-piperidin-4-yl)-1H-imidazol-2-yl]-phenyl}-3H-quinazolin-4-one;

2-{4-[1-(1-acetyl-piperidin-4-yl)-1H-imidazol-2-yl]-phenyl}-3-(5-chloro-pyridin-2-yl)-3H-quinazolin-4-one;

3-(5-chloro-pyridin-2-yl)-2-[4-(1-isopropyl-1H-imidazol-2-yl)-phenyl]-3H-quinazolin-4-one;

3-(5-chloro-pyridin-2-yl)-2-[4-(1-cyclopropyl-1H-imidazol-2-yl)-phenyl]-3H-quinazolin-4-one;

3-(5-chloro-pyridin-2-yl)-2-[4-(1-cyclobutyl-1H-imidazol-2-yl)-phenyl]-3H-quinazolin-4-one;

3-(5-chloro-pyridin-2-yl)-2-[4-(1-cyclopentyl-1H-imidazol-2-yl)-phenyl]-3H-quinazolin-4-one;

3-(5-chloro-pyridin-2-yl)-2-{4-[1-(tetrahydro-furan-3-yl)-1H-imidazol-2-yl]-phenyl}-3H-quinazolin-4-one;

3-(5-chloro-pyridin-2-yl)-2-[4-(1-pyrrolidin-3-yl-1H-imidazol-2-yl)-phenyl]-3H-quinazolin-4-one;

3-(5-chloro-pyridin-2-yl)-2-{4-[1-(1-methyl-pyrrolidin-3-yl)-1H-imidazol-2-yl]-phenyl}-3H-quinazolin-4-one;

2-{4-[1-(1-acetyl-pyrrolidin-3-yl)-1H-imidazol-2-yl]-phenyl}-3-(5-chloro-pyridin-2-yl)-3H-quinazolin-4-one;

3-(5-chloro-pyridin-2-yl)-2-{4-[1-(1-methanesulfonyl-pyrrolidin-3-yl)-1H-imidazol-2-yl]-phenyl}-3H-quinazolin-4-one;

N-[2-(2-{4-[3-(5-chloro-pyridin-2-yl)-4-oxo-3,4-dihydro-quinazolin-2-yl]-phenyl}-imidazol-1-yl)-ethyl]-acetamide;

3-(5-chloro-pyridin-2-yl)-2-{4-[1-(2-hydroxy-ethyl)-1H-imidazol-2-yl]-phenyl}-3H-quinazolin-4-one;

3-(5-chloro-pyridin-2-yl)-2-{4-[1-(2-methoxy-ethyl)-1H-imidazol-2-yl]-phenyl}-3H-quinazolin-4-one;

3-(5-chloro-pyridin-2-yl)-2-{4-[1-(2-methoxy-1-methyl-ethyl)-1H-imidazol-2-yl]-phenyl}-3H-quinazolin-4-one;

2-(2-{4-[3-(5-chloro-pyridin-2-yl)-4-oxo-3,4-dihydro-quinazolin-2-yl]-phenyl}-imidazol-1-yl)-acetamide;

2-(2-{4-[3-(5-chloro-pyridin-2-yl)-4-oxo-3,4-dihydro-quinazolin-2-yl]-phenyl}-imidazol-1-yl)-N-methyl-acetamide;

2-[4-(5-amino-furan-2-yl)-phenyl]-3-(5-chloro-pyridin-2-yl)-3H-quinazolin-4-one;

2-[4-(5-aminomethyl-furan-2-yl)-phenyl]-3-(5-chloro-pyridin-2-yl)-3H-quinazolin-4-one;

5-{4-[3-(5-chloro-pyridin-2-yl)-4-oxo-3,4-dihydro-quinazolin-2-yl]-phenyl}-furan-2-carboxylic acid amide;

2-{4-[5-(1-amino-1-methyl-ethyl)-furan-2-yl]-phenyl}-3-(5-chloro-pyridin-2-yl)-3H-quinazolin-4-one;

2-[4-(3-amino-furan-2-yl)-phenyl]-3-(5-chloro-pyridin-2-yl)-3H-quinazolin-4-one;

3-(5-chloro-pyridin-2-yl)-2-[4-(3-dimethylaminomethyl-furan-2-yl)-phenyl]-3H-quinazolin-4-one;

3-(5-chloro-pyridin-2-yl)-2-{4-[3-(1-dimethylamino-1-methyl-ethyl)-furan-2-yl]-phenyl}-3H-quinazolin-4-one;

2-{4-[3-(5-chloro-pyridin-2-yl)-4-oxo-3,4-dihydro-quinazolin-2-yl]-phenyl}-furan-3-carboxylic acid amide;

3-(5-chloro-pyridin-2-yl)-2-(4-oxazol-2-yl-phenyl)-3H-quinazolin-4-one;

2-[4-(5-aminomethyl-oxazol-2-yl)-phenyl]-3-(5-chloro-pyridin-2-yl)-3H-quinazolin-4-one;

2-{4-[3-(5-chloro-pyridin-2-yl)-4-oxo-3,4-dihydro-quinazolin-2-yl]-phenyl}-oxazole-5-carboxylic acid amide;

2-[4-(4-aminomethyl-oxazol-2-yl)-phenyl]-3-(5-chloro-pyridin-2-yl)-3H-quinazolin-4-one;

2-{4-[3-(5-chloro-pyridin-2-yl)-4-oxo-3,4-dihydro-quinazolin-2-yl]-phenyl}-oxazole-4-carboxylic acid amide;

3-(5-chloro-pyridin-2-yl)-2-(4-thiazol-2-yl-phenyl)-3H-quinazolin-4-one;

2-[4-(5-aminomethyl-thiazol-2-yl)-phenyl]-3-(5-chloro-pyridin-2-yl)-3H-quinazolin-4-one;

2-{4-[3-(5-chloro-pyridin-2-yl)-4-oxo-3,4-dihydro-quinazolin-2-yl]-phenyl}-thiazole-5-carboxylic acid amide;

2-[4-(4-amino-thiazol-2-yl)-phenyl]-3-(5-chloro-pyridin-2-yl)-3H-quinazolin-4-one;

N-(2-{4-[3-(5-chloro-pyridin-2-yl)-4-oxo-3,4-dihydro-quinazolin-2-yl]-phenyl}-thiazol-4-yl)-acetamide;

2-[4-(5-amino-thiazol-2-yl)-phenyl]-3-(5-chloro-pyridin-2-yl)-3H-quinazolin-4-one;

N-(2-{4-[3-(5-chloro-pyridin-2-yl)-4-oxo-3,4-dihydro-quinazolin-2-yl]-phenyl}-thiazol-5-yl)-acetamide;

3-(5-chloro-pyridin-2-yl)-2-[4-(2-oxo-tetrahydro-pyrimidin-1-yl)-phenyl]-3H-quinazolin-4-one;

3-(5-chloro-pyridin-2-yl)-2-[4-(2-oxo-imidazolidin-1-yl)-phenyl]-3H-quinazolin-4-one;

3-(5-chloro-pyridin-2-yl)-2-[4-(2-oxo-[1,3]diazepan-1-yl)-phenyl]-3H-quinazolin-4-one;

2-[4-(3-amino-2-oxo-piperidin-1-yl)-phenyl]-3-(5-chloro-pyridin-2-yl)-3H-quinazolin-4-one;

3-(5-chloro-pyridin-2-yl)-2-[4-(3-dimethylamino-2-oxo-piperidin-1-yl)-phenyl]-3H-quinazolin-4-one;

3-(5-chloro-pyridin-2-yl)-2-[4-(2-oxo-3-pyrrolidin-1-yl-piperidin-1-yl)-phenyl]-3H-quinazolin-4-one;

N-(1-{4-[3-(5-chloro-pyridin-2-yl)-4-oxo-3,4-dihydro-quinazolin-2-yl]-phenyl}-2-oxo-piperidin-3-yl)-acetamide;

2-[4-(3-amino-2-oxo-pyrrolidin-1-yl)-phenyl]-3-(5-chloro-pyridin-2-yl)-3H-quinazolin-4-one;

3-(5-chloro-pyridin-2-yl)-2-[4-(2-dimethylaminomethyl-imidazol-1-yl)-phenyl]-3H-quinazolin-4-one;

1-{4-[3-(5-chloro-pyridin-2-yl)-4-oxo-3,4-dihydro-quinazolin-2-yl]-phenyl}-1H-imidazole-2-carboxylic acid dimethylamide;

3-(5-chloro-pyridin-2-yl)-2-(4-isoxazol-5-yl-phenyl)-3H-quinazolin-4-one;

3-(5-chloro-pyridin-2-yl)-2-(4-oxazol-5-yl-phenyl)-3H-quinazolin-4-one;

3-(5-chloro-pyridin-2-yl)-2-(4-thiazol-5-yl-phenyl)-3H-quinazolin-4-one;

3-(5-chloro-pyridin-2-yl)-2-[4-(3H-[1,2,3]triazol-4-yl)-phenyl]-3H-quinazolin-4-one;

3-(5-chloro-pyridin-2-yl)-2-[4-(5-methyl-4H-[1,2,4]triazol-3-yl)-phenyl]-3H-quinazolin-4-one;

3-(5-chloro-pyridin-2-yl)-2-[4-(5-methyl-[1,3,4]thiadiazol-2-yl)-phenyl]-3H-quinazolin-4-one;

2-biphenyl-4-yl-3-(5-chloro-pyridin-2-yl)-3H-quinazolin-4-one;

2-(2'-amino-biphenyl-4-yl)-3-(5-chloro-pyridin-2-yl)-3H-quinazolin-4-one;

3-(5-chloro-pyridin-2-yl)-2-(2'-dimethylamino-biphenyl-4-yl)-3H-quinazolin-4-one;

2-(2'-aminomethyl-biphenyl-4-yl)-3-(5-chloro-pyridin-2-yl)-3H-quinazolin-4-one;

3-(5-chloro-pyridin-2-yl)-2-(2'-dimethylaminomethyl-biphenyl-4-yl)-3H-quinazolin-4-one;

3-(5-chloro-pyridin-2-yl)-2-(4-pyridin-2-yl-phenyl)-3H-quinazolin-4-one;

3-(5-chloro-pyridin-2-yl)-2-(4-pyridin-3-yl-phenyl)-3H-quinazolin-4-one;

3-(5-chloro-pyridin-2-yl)-2-(4-pyridin-4-yl-phenyl)-3H-quinazolin-4-one;

2-[4-(2-amino-pyridin-3-yl)-phenyl]-3-(5-chloro-pyridin-2-yl)-3H-quinazolin-4-one;

2-[4-(2-aminomethyl-pyridin-3-yl)-phenyl]-3-(5-chloro-pyridin-2-yl)-3H-quinazolin-4-one;

3-(5-chloro-pyridin-2-yl)-2-[4-(2-dimethylaminomethyl-pyridin-3-yl)-phenyl]-3H-quinazolin-4-one;

3-(5-chloro-pyridin-2-yl)-2-[2-fluoro-4-(2-oxo-2H-pyridin-1-yl)-phenyl]-3H-quinazolin-4-one;

3-(5-chloro-pyridin-2-yl)-2-[2-fluoro-4-(2-oxo-piperidin-1-yl)-phenyl]-3H-quinazolin-4-one;

6-chloro-3-(5-chloro-pyridin-2-yl)-2-[2-fluoro-4-(2-oxo-piperidin-1-yl)-phenyl]-3H-quinazolin-4-one;

3-(5-chloro-pyridin-2-yl)-6-fluoro-2-[2-fluoro-4-(2-oxo-piperidin-1-yl)-phenyl]-3H-quinazolin-4-one;

6-bromo-3-(5-chloro-pyridin-2-yl)-2-[2-fluoro-4-(2-oxo-piperidin-1-yl)-phenyl]-3H-quinazolin-4-one;

3-(5-chloro-pyridin-2-yl)-2-[2-fluoro-4-(2-oxo-piperidin-1-yl)-phenyl]-4-oxo-3,4-dihydro-quinazoline-6-carbonitrile;

3-(5-chloro-pyridin-2-yl)-2-[2-fluoro-4-(2-oxo-piperidin-1-yl)-phenyl]-6-methoxy-3H-quinazolin-4-one;

3-(5-chloro-pyridin-2-yl)-2-[2-fluoro-4-(2-oxo-piperidin-1-yl)-phenyl]-4-oxo-3,4-dihydro-quinazoline-6-carboxylic acid amide;

6-chloro-3-(5-chloro-pyridin-2-yl)-2-[2-fluoro-4-(2-oxo-2H-pyridin-1-yl)-phenyl]-3H-quinazolin-4-one;

3-(5-chloro-pyridin-2-yl)-2-[2-fluoro-4-(2-oxo-2H-pyridin-1-yl)-phenyl]-6-methoxy-3H-quinazolin-4-one;

3-(5-chloro-pyridin-2-yl)-6-fluoro-2-[2-fluoro-4-(2-oxo-2H-pyridin-1-yl)-phenyl]-3H-quinazolin-4-one;

3-(5-chloro-pyridin-2-yl)-2-[2-fluoro-4-(2-oxo-2H-pyridin-1-yl)-phenyl]-4-oxo-3,4-dihydro-quinazoline-6-carbonitrile;

3-(4-chloro-phenyl)-2-[2-fluoro-4-(2-oxo-2H-pyridin-1-yl)-phenyl]-3H-quinazolin-4-one;

2-[2-fluoro-4-(2-oxo-2H-pyridin-1-yl)-phenyl]-3-(4-methoxy-phenyl)-3H-quinazolin-4-one;

3-(3-chloro-phenyl)-2-[2-fluoro-4-(2-oxo-2H-pyridin-1-yl)-phenyl]-3H-quinazolin-4-one;

2-fluoro-5-{2-[2-fluoro-4-(2-oxo-2H-pyridin-1-yl)-phenyl]-4-oxo-4H-quinazolin-3-yl}-benzonitrile;

3-(5-chloro-pyridin-2-yl)-2-[4-(2-oxo-pyrrolidin-1-yl)-phenyl]-3H-quinazolin-4-one;

3-(4-chloro-phenyl)-2-[4-(1H-pyrrol-2-yl)-phenyl]-3H-quinazolin-4-one;

3-(4-methoxy-phenyl)-2-[4-(1H-pyrrol-2-yl)-phenyl]-3H-quinazolin-4-one;

3-(3-chloro-phenyl)-2-[4-(1H-pyrrol-2-yl)-phenyl]-3H-quinazolin-4-one;

3-(3-chloro-phenyl)-2-[4-(1-methyl-1H-pyrrol-2-yl)-phenyl]-3H-quinazolin-4-one;

3-(4-chloro-phenyl)-2-[4-(1-methyl-1H-pyrrol-2-yl)-phenyl]-3H-quinazolin-4-one;

3-(5-chloro-pyridin-2-yl)-6-methoxy-2-[4-(1-methyl-1H-pyrrol-2-yl)-phenyl]-3H-quinazolin-4-one;

3-(5-chloro-pyridin-2-yl)-2-{4-[1-(2-dimethylamino-ethyl)-1H-pyrrol-2-yl]-phenyl}-6-methoxy-3H-quinazolin-4-one;

6-chloro-3-(5-chloro-pyridin-2-yl)-2-{4-[1-(2-dimethylamino-ethyl)-1H-pyrrol-2-yl]-phenyl}-3H-quinazolin-4-one;

6-chloro-3-(4-chloro-phenyl)-2-{4-[1-(2-dimethylamino-ethyl)-1H-pyrrol-2-yl]-phenyl}-3H-quinazolin-4-one;

3-(5-chloro-pyridin-2-yl)-6-methoxy-2-[4-(2-oxo-imidazolidin-1-yl)-phenyl]-3H-quinazolin-4-one;

6-chloro-3-(5-chloro-pyridin-2-yl)-2-[4-(2-oxo-imidazolidin-1-yl)-phenyl]-3H-quinazolin-4-one;

6-chloro-3-(4-chloro-phenyl)-2-[4-(2-oxo-imidazolidin-1-yl)-phenyl]-3H-quinazolin-4-one;

6-chloro-3-(4-methoxy-phenyl)-2-[4-(2-oxo-imidazolidin-1-yl)-phenyl]-3H-quinazolin-4-one;

6-chloro-3-(5-chloro-pyridin-2-yl)-2-[4-(2-oxo-tetrahydro-pyrimidin-1-yl)-phenyl]-3H-quinazolin-4-one;

3-(5-chloro-pyridin-2-yl)-6-methoxy-2-[4-(2-oxo-tetrahydro-pyrimidin-1-yl)-phenyl]-3H-quinazolin-4-one;

3-(5-chloro-pyridin-2-yl)-2-[4-(2-dimethylaminomethyl-4,5-dihydro-imidazol-1-yl)-phenyl]-6-methoxy-3H-quinazolin-4-one;

3-(4-chloro-phenyl)-2-[4-(2-dimethylaminomethyl-4,5-dihydro-imidazol-1-yl)-phenyl]-6-methoxy-3H-quinazolin-4-one;

6-chloro-3-(4-chloro-phenyl)-2-[4-(2-dimethylaminomethyl-4,5-dihydro-imidazol-1-yl)-phenyl]-3H-quinazolin-4-one;

6-bromo-3-(4-chloro-phenyl)-2-[4-(2-dimethylaminomethyl-4,5-dihydro-imidazol-1-yl)-phenyl]-3H-quinazolin-4-one;

2-[4-(2-dimethylaminomethyl-4,5-dihydro-imidazol-1-yl)-phenyl]-3-(4-methoxy-phenyl)-6-methyl-3H-quinazolin-4-one;

3-(5-chloro-pyridin-2-yl)-2-[4-(1H-pyrrol-2-yl)-phenyl]-3H-pyrido[3,2-d]pyrimidin-4-one;

3-(5-chloro-pyridin-2-yl)-2-[4-(1-methyl-1H-pyrrol-2-yl)-phenyl]-3H-pyrido[3,2-d]pyrimidin-4-one;

3-(5-chloro-pyridin-2-yl)-2-{4-[1-(2-methylamino-ethyl)-1H-pyrrol-2-yl]-phenyl}-3H-pyrido[3,2-d]pyrimidin-4-one;

3-(5-chloro-pyridin-2-yl)-2-[4-(1H-pyrrol-2-yl)-phenyl]-3H-pyrido[4,3-d]pyrimidin-4-one;

3-(5-chloro-pyridin-2-yl)-2-[4-(1-methyl-1H-pyrrol-2-yl)-phenyl]-3H-pyrido[4,3-d]pyrimidin-4-one;

3-(5-chloro-pyridin-2-yl)-2-[4-(1H-pyrrol-2-yl)-phenyl]-3H-pyrido[3,4-d]pyrimidin-4-one;

3-(5-chloro-pyridin-2-yl)-2-[4-(1H-pyrrol-2-yl)-phenyl]-3H-thieno[3,4-d]pyrimidin-4-one;

3-(5-chloro-pyridin-2-yl)-2-[4-(1H-imidazol-2-yl)-phenyl]-3H-pyrido[3,2-d]pyrimidin-4-one;

3-(5-chloro-pyridin-2-yl)-2-[4-(1H-imidazol-2-yl)-phenyl]-3H-thieno[3,4-d]pyrimidin-4-one;

3-(5-chloro-pyridin-2-yl)-2-[4-(1-methyl-1H-imidazol-2-yl)-phenyl]-3H-thieno[3,4-d]pyrimidin-4-one;

3-(5-chloro-pyridin-2-yl)-2-[4-(2-oxo-tetrahydro-pyrimidin-1-yl)-phenyl]-3H-pyrido[3,2-d]pyrimidin-4-one;

3-(5-chloro-pyridin-2-yl)-2-[4-(2-oxo-tetrahydro-pyrimidin-1-yl)-phenyl]-3H-pyrido[4,3-d]pyrimidin-4-one;

3-(5-chloro-pyridin-2-yl)-2-[4-(2-oxo-tetrahydro-pyrimidin-1-yl)-phenyl]-3H-pyrido[3,4-d]pyrimidin-4-one;

3-(5-chloro-pyridin-2-yl)-2-[4-(2-oxo-tetrahydro-pyrimidin-1-yl)-phenyl]-3H-thieno[3,4-d]pyrimidin-4-one;

3-(5-chloro-pyridin-2-yl)-2-[4-(2-oxo-imidazolidin-1-yl)-phenyl]-3H-pyrido[3,2-d]pyrimidin-4-one;

3-(5-chloro-pyridin-2-yl)-2-[4-(2-oxo-imidazolidin-1-yl)-phenyl]-3H-pyrido[4,3-d]pyrimidin-4-one;

6-(5-chloro-pyridin-2-yl)-5-[4-(2-oxo-tetrahydro-pyrimidin-1-yl)-phenyl]-6H-[1,2,5]oxadiazolo[3,4-d]pyrimidin-7-one;

6-(5-chloro-pyridin-2-yl)-5-[4-(2-oxo-tetrahydro-pyrimidin-1-yl)-phenyl]-6H-isoxazolo[4,3-d]pyrimidin-7-one;

6-(5-chloro-pyridin-2-yl)-5-[4-(2-oxo-tetrahydro-pyrimidin-1-yl)-phenyl]-6H-thiazolo[5,4-d]pyrimidin-7-one;

1-(5-chloro-pyridin-2-yl)-7-methyl-2-[4-(2-oxo-tetrahydro-pyrimidin-1-yl)-phenyl]-1,7-dihydro-purin-6-one;

3-(5-chloro-pyridin-2-yl)-2-[4-(2-oxo-imidazolidin-1-yl)-phenyl]-3H-thieno[3,2-d]pyrimidin-4-one;

3-(5-chloro-pyridin-2-yl)-5-methyl-2-[4-(2-oxo-imidazolidin-1-yl)-phenyl]-3,5-dihydro-pyrrolo[3,2-d]pyrimidin-4-one;

6-(5-chloro-pyridin-2-yl)-1-methyl-5-[4-(2-oxo-imidazolidin-1-yl)-phenyl]-1,6-dihydro-pyrazolo[4,3-d]pyrimidin-7-one;

3-(5-chloro-pyridin-2-yl)-2-[4-(3-dimethylamino-2-oxo-piperidin-1-yl)-phenyl]-3H-furo[3,2-d]pyrimidin-4-one;

3-(5-chloro-pyridin-2-yl)-2-[4-(3-dimethylamino-2-oxo-piperidin-1-yl)-phenyl]-5-methyl-3,5-dihydro-pyrrolo[3,2-d]pyrimidin-4-one; and,

6-(5-chloro-pyridin-2-yl)-5-[4-(3-dimethylamino-2-oxo-piperidin-1-yl)-phenyl]-1-methyl-1,6-dihydro-pyrazolo[4,3-d]pyrimidin-7-one;

or a pharmaceutically acceptable salt form thereof.

9. A method for treating a thromboembolic disorder, comprising: administering to a patient in need thereof a therapeutically effective amount of a compound of Claim 1or a pharmaceutically acceptable salt thereof.

10. A method according to Claim 9, wherein the thromboembolic disorder is selected from the group consisting of arterial cardiovascular thromboembolic disorders, venous cardiovascular thromboembolic disorders, and thromboembolic disorders in the chambers of the heart.

11. A method according to Claim 9, wherein the thromboembolic disorder is selected from unstable angina, an acute coronary syndrome, first myocardial infarction, recurrent myocardial infarction, ischemic sudden death, transient ischemic attack, stroke, atherosclerosis, peripheral occlusive arterial disease, venous thrombosis, deep vein thrombosis, thrombophlebitis, arterial embolism, coronary arterial thrombosis, cerebral arterial thrombosis, cerebral embolism, kidney embolism, pulmonary embolism, and thrombosis resulting from (a) prosthetic valves or other implants, (b) indwelling catheters, (c) stents, (d) cardiopulmonary bypass, (e) hemodialysis, and (f) other procedures in which blood is exposed to an artificial surface that promotes thrombosis.

12. A method of treating a patient in need of thromboembolic disorder treatment, comprising: administering a compound of Claim 1or a pharmaceutically acceptable salt thereof in an amount effective to treat a thromboembolic disorder.

13. A method, comprising: administering a compound of Claim 1or a pharmaceutically acceptable salt thereof in an amount effective to treat a thromboembolic disorder.